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SOME RESULTS IN THE THEORY OF SUBSET SELECTION PROCEDURES, (U)

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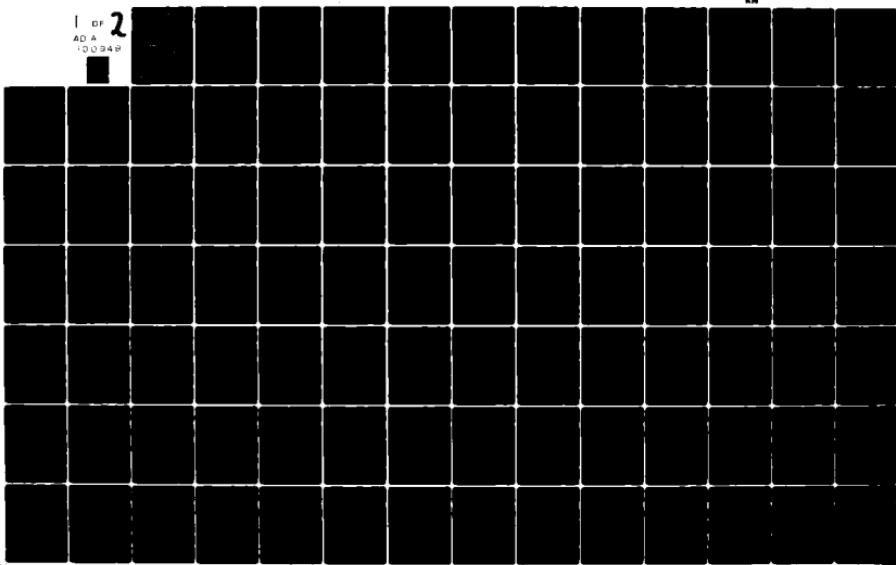
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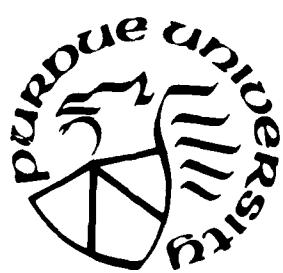


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Some Results in the Theory
of Subset Selection Procedures

by

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INTRODUCTION

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Selection and ranking (ordering) problems in statistical inference arise mainly because the classical tests of homogeneity are often inadequate in certain situations where the experimenter is interested in comparing $k \geq 2$ populations, treatments or processes with the goal of selecting one or more worthwhile (good) populations. Mosteller (1949), Paulson (1949), Bahadur (1950) and Bahadur and Robbins (1950) were among the earliest research workers to recognize this inadequacy and to formulate the problem as a multiple decision problem aimed at the selection and ranking of the k populations.

In the thirty years since these early papers, selection and ranking problems have become an active area of statistical research. There have been two approaches to these problems, the 'indifference zone' approach and the 'subset selection' approach. In the first approach, due to Bechhofer (1954), the experimenter wishes to select one population (or a fixed number $t \geq 1$ of population) which is guaranteed to be the one of interest to him with a fixed probability P^* whenever the unknown parameters lie outside some subspace of the parameter space, the so-called indifference zone. Important contributions using this approach have been made by Bechhofer and Sobel (1954), Bechhofer, Dunnett and Sobel (1954), Sobel and Huyett (1957), Sobel (1962), Bechhofer, Trotter and Sobel (1968), Mahanunulu (1967), Desu and Sobel (1962, 1971) and

Tamhane and Bechhofer (1977, 1979) among others. A quite complete bibliography may be found in Gupta and Panchapakesan (1979) (see also Gibbons, Olkin and Sobel (1977)).

The second approach pioneered by Gupta (1956, 1963, 1965) assumes no a priori information about the parameter space. A single population is not necessarily chosen; rather a subset of the given k populations is selected depending on the outcome of the experiment. It is guaranteed to contain the population(s) of interest with probability which is at least equal to P^* (the basic probability requirement) regardless of the true unknown configurations of the parameters. Some recent contributors in the category of subset selection include: Beely (1965), Gnanadesikan (1966), Gnanadesikan and Gupta (1970), Gupta (1967), Gupta and Studden (1970), Nagel (1970), Gupta and Nagel (1971), Gupta and Panchapakesan (1972), Rizvi and Sobel (1967), McDonald (1969), Gupta and McDonald (1970), Santner (1975), W. T. Huang (1972), D. Y. Huang (1975), Gupta and Huang (1975a, 1975b) and Gupta and Huang (1976).

Subset selection procedures can also be thought of as screening procedures which enable the experimenter to select a subset of populations (under study) which contains the populations of interest so that the populations in the selected subset can be further studies.

Sequential and multistage aspects of the ranking and selection problems, have been explored, based on the indifference zone approach by Bechhofer, Dunnett and Sobel (1954), Bechhofer (1958), Paulson (1962, 1963, 1964, 1967) and Bechhofer, Fiefer and Sobel (1968). Barson and Gupta (1972), Huang (1972), Gupta and Huang (1975), Gupta and Miescke (1979) and Carroll (1974) have investigated subset selection procedures, based on sequential sampling.

Contributions to optimum properties of subset selection procedures have been made by Lehmann (1961), Studden (1967), Deely and Gupta (1968), Berger (1977, 1979), Gupta and Hsu (1978), Gupta and Miescke (1978), Berger and Gupta (1980).

In the decision-theoretic approach to the subset selection problem, Goel and Rubin (1977), Chernoff and Yahav (1977), Bickel and Yahav (1977), Gupta and Hsu (1978), Miescke (1979), Gupta and Kim (1980), Gupta and Hsiao (1980) have given different formulations under different loss functions and carried out investigations which indicate that the gamma type maximum (minimum) means procedures are quite "optimal" and "robust".

The main purpose of this thesis is to study some problems arising in the subset selection approach and provide procedures and results for some unsolved problems.

Chapter I considers the problem of selecting a subset containing all populations better than a control under an ordering prior. Here, by an ordering prior we mean that there exists a known simple or partial order relationship among the unknown parameters of the treatments (excluding the control). Three new selection procedures are proposed and studied. These procedures do meet the usual requirement that the probability of a correct selection is greater than or equal to a pre-determined number P^* . Two of the three procedures use the isotonic regression over the sample means of the k treatments with respect to (wrt) the given ordering prior. Tables which are necessary to carry out the selection procedures with isotonic approach for the selection of unknown means of normal populations and gamma populations are given. Monte Carlo comparisons on the performance of several procedures for the normal or gamma means problem were carried out in

several selected cases; these are given in Table V and Table VI at the end of Chapter I. In each case ten thousand simulations were performed. The results of this study seem to indicate that the procedures based on isotonic estimators always have superior performance, especially, when there are more than one bad populations (in comparison with the control).

Chapter II deals with a new 'Bayes-P*' approach about the problem of selecting a subset which contains the 'best' of k populations. Here, by best we mean the (unknown) population with the largest unknown mean. The (non-randomized) Bayes-P* rule refers to a rule with minimum risk in the class of (non-randomized) rules which satisfy the condition that the posterior probability of selecting the best is at least equal to P^* . Given the priors of the unknown parameters, two 'Bayes-P*' subset selection procedures ψ^B and ψ_{NR}^B (randomized and non-randomized, respectively) under certain loss functions are obtained and compared with the classical maximum-type means procedure ψ^M . The comparisons of the performance of ψ^B with ψ_{NR}^B and ψ^M , based on Monte Carlo studies, indicate that the procedure ψ^B always has higher 'efficiency' and smaller expected selected size of the selected subset. Also ψ^B appears to be robust when the true distributions are not normal but are some other symmetric distributions such as, the logistic, the double exponential, Laplace, and the gross error model (the contaminated distribution).

CHAPTER I
SELECTION PROCEDURES FOR POPULATIONS
BETTER THAN A CONTROL UNDER ORDERING PRIOR

1.1. Introduction

In this chapter, three new selection procedures are given for the problem of selecting a subset which contains all populations better than a standard or control under simple or partial ordering prior. Here by simple or partial ordering prior we mean that there exist known simple or partial order relationships (defined more specifically later in Section 1.2) among unknown parameters. The procedures described do meet the usual requirement that the probabilities of a correct selection are greater than or equal to a predetermined number P^* , the so-called P^* condition.

Many authors have considered the problem of comparing populations with a control under different types of formulations (see Gupta and Panchapakesan (1979)). Dunnett (1955) considered the problem of separating those treatments which are better than the control from those that are worse. Gupta and Sobel (1958), Gupta (1965), Naik (1975), Broström (1977) studied the problem of selecting a subset containing all populations better than the control. Lehmann (1961) discussed similar problems with emphasis on the derivation of a restricted minimax procedure. Kim (1979), Hsiao (1979) studied the problem of

selecting populations close to a control. In all these papers it is assumed that all populations are independent and that there is no information about the order of unknown parameters. However, in many situations, we may know something about the unknown parameters. What we know is always not the prior distributions but some partial or incomplete prior information, such as the simple or partial order relationship among the unknown parameters. This type of information about the ordering prior may come from the past experiences; or it may arise in the experiments where, for example, higher dose level of some drugs always has larger effect (side-effect) on the patients.

In Section 1.2 definitions and notations used in this chapter are introduced. In Section 1.3 we consider the problem for location parameters. We propose three types of selection procedures for the cases when the control parameter is known or not known (the scale parameter may or may not be assumed known). Some equivalent forms of the procedures are given, and their properties are discussed. In Section 1.4 the problem for scale parameters of the gamma distributions is considered and three analogous selection procedures are proposed. In both Section 1.3 and 1.4 simple ordering priors are assumed and some theorems in the theory of random walks are used. In Section 1.5 a selection procedure is given for the problem of selecting all populations better than the control under partial ordering prior. Section 1.6 deals with the use of Monte Carlo techniques to make comparisons among the selection procedures proposed in Section 1.3 and those in Section 1.4, respectively.

1.2. Notations and Definitions

Suppose we have $k + 1$ populations " $\pi_0, \pi_1, \dots, \pi_k$ ". The population treatment " π_0 " is called the control or standard population. Assume that the random variables X_{ij} associated with $F(\cdot; \pi_i)$ and x_{i1}, \dots, x_{in_i} , $i = 1, \dots, k$, is an independent sample from π_i . Assume that we have an ordering prior of " π_1, \dots, π_k ". First we assume that the ordering prior is the simple order, so that without loss of generality, we may assume that, $\pi_1 < \dots < \pi_k$. In Section 1.5 we will consider the partial ordering prior case. Note that the values of π_i 's are unknown.

Suppose our goal is to find a non-trivial (small) subset which contains all populations with parameter larger (smaller) than the control " π_0 " (known or unknown) with probability not less than a given value α^* .

The action space \mathcal{C} is the class of all subsets of set $\{1, 2, \dots, k\}$. An action A is the selection of some subset of the k populations. This means that π_j is included in the selected subset.

Let $\pi = (\pi_0, \pi_1, \dots, \pi_k)$. Then the parameter space is denoted by Σ , where $\Sigma = \{\pi \in \mathbb{R}^{k+1} | \pi_1 < \pi_2 < \dots < \pi_k; -\infty < \pi_0 < \infty\}$ is a subset of $k + 1$ dimensional Euclidean space \mathbb{R}^{k+1} .

The sample space is denoted by Ω where

$$\Omega = \{x \in \mathbb{R}^{n_1 + \dots + n_k} | x = (x_{11}, \dots, x_{1n_1}, x_{21}, \dots, x_{kn_k}, \dots, x_{kn_k})\}.$$

Definition 1.2.1. A (non-randomized) selection procedure (rule) δ is a mapping from Ω to \mathcal{C} .

A population " π_i " ($i = 1, \dots, k$) is called a good population if " $\pi_i \geq \pi_0$ ", and we say a selection procedure δ make a correct selection (CS) if the selected subset contains all good populations. A selection procedure δ satisfies the P^* -condition if

$$P_n(\text{CS}|\delta) \geq P^* \quad \text{for all } \theta \in \Theta$$

that is

$$\inf_{\theta \in \Theta} P_n(\text{CS}|\delta) \geq P^*. \quad (1.2.1)$$

Let $\Delta = \{\delta \mid \inf_{\theta \in \Theta} P_n(\text{CS}|\delta) \geq P^*\}$ be a collection of all selection procedures satisfying the P^* -condition.

In the sequel we will use the isotonic estimators (see Barlow, Bartholomew, Bremner and Brunk (1972)). Hence we give the following definitions and theorems.

Definition 1.2.2. Let the set \mathcal{S} be a finite set. A binary relation " \leq " on \mathcal{S} is called a simple order if it is

- (1) reflexive: $x \leq x$ for $x \in \mathcal{S}$
- (2) transitive: $x, y, z \in \mathcal{S}$ and $x \leq y, y \leq z$ imply $x \leq z$
- (3) antisymmetric: $x, y \in \mathcal{S}$ and $x \leq y, y \leq x$ imply $x = y$
- (4) every two elements are comparable: $x, y \in \mathcal{S}$ imply either $x \leq y$ or $y \leq x$.

A partial order on \mathcal{S} is a binary relation " \leq " on \mathcal{S} , such that it is (1) reflexive, (2) transitive, and (3) antisymmetric. Thus every simple order is a partial order. We use poset (\mathcal{S}, \leq) to denote the set \mathcal{S} that has a partial order binary relation " \leq " on it.

Definition 1.2.3. A real-valued function f is called isotonic on poset (\mathcal{S}, \leq) if and only if (1) f is defined on \mathcal{S} , (2) if $x, y \in \mathcal{S}$ and $x \leq y$ then $f(x) \leq f(y)$.

Definition 1.2.4. Let g be a real-valued function on \mathcal{S} and let w be a given positive function on \mathcal{S} . A function g^* on \mathcal{S} is called an isotonic regression of g with weights w if and only if:

(1) g^* is an isotonic function on poset (\mathcal{S}, \leq)

$$(2) \quad \int_{\mathcal{S}} [g(x) - g^*(x)]^2 w(x) = \min_{f \in \mathcal{G}(\mathcal{S})} \int_{\mathcal{S}} [g(x) - f(x)]^2 w(x),$$

where \mathcal{G} is the class of all isotonic functions on poset (\mathcal{S}, \leq) .

From Barlow, et. al. (1972), (see their Theorem 1, Corollary and the corollary there), we have the following theorem.

Theorem 1.2.1. There exists one and only one isotonic regression g^* of g with weight w on poset (\mathcal{S}, \leq) .

Definition 1.2.5. A set S is convex if $s_1, s_2 \in S$ and if $0 < t < 1$ then $s_1 + (1-t)s_2 \in S$.

Definition 1.2.6. A set S is a cone if $s \in S$ then for any non-negative real number c , $cs \in S$.

Definition 1.2.7. A poset (\mathcal{S}, \leq) is a lattice if for all A and B in \mathcal{S} there exist for any finite non-empty subset H of \mathcal{S} ,

If f and t are two isotonic functions on poset (\mathcal{S}, \leq) , we define $f \wedge q$ and $t \vee q$ as

$$(f \wedge q)(t) = f(t) \wedge q(t) = \min(f(t), q(t))$$

and

$$(f \vee g)(t) = r(t) \vee g(t) = \max(r(t), g(t)).$$

Then we state the following:

Theorem 1.2.2. The class δ of all isotonic functions on poset (\mathbb{S}, \leq) is a convex cone and a lattice.

There are some algorithms, such as the "pool-adjacent-violation" algorithm (see page 13 of Barlow, et. al. (1972)) or Ayer, Frank, Tukey, Reid and Silverman (1955) or the "up-and-down blocks" algorithm, see Fager (1964), which show how to calculate the isotonic regression under strict order.

The following max-min formulas were given by Ayer et. al. (1955).

Theorem 1.2.3. (max-min formulas)

Assume that we have poset (\mathbb{S}, \leq) where $\mathbb{S} = \{1, \dots, k\} \cup \{\infty\}$, and that function $g: \mathbb{S} \rightarrow \mathbb{R}$, then the isotonic regression g^* of g with weight W has the following formulas:

$$\begin{aligned} g^*(i) &= \max_{s \leq i \leq t} \text{Av}(s, t) \\ &= \max_{s \leq i \leq t} \min_{s \leq s' \leq t} \text{Av}(s, t) \\ &= \min_{t \geq i} \max_{s \leq s' \leq t} \text{Av}(s, t) \\ &= \min_{t \geq i} \max_{s \leq s' \leq t} \text{Av}(s, t) \end{aligned}$$

where

$$\text{Av}(s, t) = \frac{\sum_{r=s}^t w(r)}{\sum_{r=s}^t w(r)},$$

$$\begin{cases} \sum_{r=s}^t w(r) \\ \sum_{r=s}^t w(r) \end{cases}$$

Corollary 1.2.1. $(q + c)^* \leq q^* + c$,

$$(aq)^* = aq^*, \text{ if } a > 0.$$

Corollary 1.2.2. $[(\eta^*)q + \zeta(\eta^*)]^* \leq (\eta^*)q^* + \zeta(\eta^*)$, where η^* is a non-negative function and ζ is an arbitrary function.

1.3. Proposed Selection Procedures for the Location Parameter Problem

To discuss some more general results, we assume that population π_j has an absolutely continuous location-scale distribution function

$F(x; \omega_1, \dots, \omega_k) = F(\frac{x-\omega_1}{\omega_2}, \dots, \frac{\omega_{k-1}-\omega_k}{\omega_k})$, where $0 < F(x) < 1$ for all x , $\omega_1 > \omega_2 > \dots > \omega_k$ and the values of $\omega_1, \dots, \omega_k$ are unknown, but their ordering, say, $\omega_1 > \dots > \omega_k$ is known. Note that in this case we replace ω in the parameter space by ω , all other quantities remaining the same.

Let us define the subspace $\pi_i = \{\omega \in \mathbb{R}^k : \omega_{k-i+1} = \omega_0 = \text{fixed}\}$ for $i = 1, \dots, k-1$ and let subspace $\pi_k = \{\omega \in \mathbb{R}^k : \omega_0 = \omega_1\}$ and suppose

$\omega_0 = \omega \in \pi_k \cap \pi_0$, then we have $\omega = \sum_{i=0}^{k-1} \omega_i e_i$. Note that the constant ω_0 could be known or unknown. If ω_0 is unknown, we assume that the distribution of population π_0 is $F(x; \omega_0, \sigma^2)$ and we take independent observations x_{01}, \dots, x_{0n_0} from π_0 and the sample space Ω turns to

$\{x \in \mathbb{R}^{n_0+\dots+n_k} | x = (x_{01}, \dots, x_{0n_0}, x_{11}, \dots, x_{kn_k})\}$. Using the partition π_0, \dots, π_k of parameter space π , we have

$$\inf_{\pi \in \Omega} P_\pi(CS^*) = \inf_{\pi \in \Omega} \inf_{1 \leq i \leq k} P_\pi(CS^*_i),$$

for any selection procedure CS^* . Hence the P^* -condition is equivalent to

$$\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|s) \geq P^*, \text{ for } i = 1, \dots, k.$$

Note that $\inf_{\mu \in \mathcal{C}_0} P_\mu(CS|s) = 1$ for any selection procedure since there

exist no good population in this case.

Suppose $X_i = x_i$ is the outcome of the sample mean of population μ_i , $i = 1, \dots, k$. Let s denote the set $\{\mu_1, \mu_2, \dots, \mu_k\}$ where $\mu_1 < \dots < \mu_k$, and let $w(\mu_i) = n_i^{-1/2} - w_i$, $g(\mu_i) = x_i$, $i = 1, \dots, k$. Then by the max-min formulas, the isotonic regression of g is g^* , where

$$g^*(\mu_i) = \max_{1 \leq s \leq i} \min_{s \leq t \leq k} \frac{\sum_{j=s}^t x_j w_j}{\sum_{j=s}^t w_j}, \quad i = 1, \dots, k.$$

The isotonic estimator of μ_i is denoted by $\hat{\mu}_{i:k}$, $i = 1, \dots, k$ where

$$\begin{aligned} \hat{\mu}_{i:k} &= \max_{1 \leq s \leq i} \min_{s \leq t \leq k} \frac{\sum_{j=s}^t x_j w_j}{\sum_{j=s}^t w_j} \\ &= \max_{1 \leq j \leq i} \hat{x}_{j:k}. \end{aligned} \quad (1.3.1)$$

where

$$\hat{x}_{j:k} = \min \left(x_j, \frac{x_j w_j + x_{j+1} w_{j+1}}{w_j + w_{j+1}}, \dots, \frac{x_j w_j + \dots + x_k w_k}{w_j + \dots + w_k} \right). \quad (1.3.2)$$

1.3.1. Proposed Selection Procedure

Case I. σ_0^2 known, common variance σ^2 known, and common sample size n .

Definition 1.3.1. We define the procedure π_1 as follows:

Step 1. Select π_i , $i = 1, \dots, k$ and stop, if

$$\hat{x}_{1:k} > \pi_0 - d_{1:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject π_1 and go to step 2.

Step 2. Select π_i , $i = 2, \dots, k$ and stop, if

$$\hat{x}_{2:k} > \pi_0 - d_{2:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject π_2 and go to step 3.

:

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Step $k-1$. Select π_i , $i = k-1, k$ and stop, if

$$\hat{x}_{k-1:k} > \pi_0 - d_{k-1:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject π_{k-1} and go to step k .

Step k . Select π_k and stop, if

$$\hat{x}_{k:k} > \pi_0 - d_{k:k}^{(1)} \frac{\sigma}{\sqrt{n}},$$

otherwise reject π_k .

Here $d_{i:k}^{(1)}$'s are the smallest values such that $\pi_1 \in \Delta$, that is, π_1 satisfies the p^* -condition.

1.3.2. On the Evaluation of $\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i+1})$ and the Value of the

Constants $d_{1:k}^{(1)}, \dots, d_{k:k}^{(1)}$

For any $\mu \in \mathcal{C}_i$, $1 \leq i \leq k$, let Z_i 's i.i.d. $F(\cdot; 0, 1)$ then

$$\begin{aligned} & P_\mu(CS|_{i+1}) \\ &= P_\mu\left(\bigcup_{j=1}^{k-i+1} \{X_{j:k} \geq \mu_0 - d_{j:k}^{(1)} \frac{\sigma}{\sqrt{n}}\}\right) \\ &= P_\mu\left(\bigcup_{j=1}^{k-i+1} \bigcup_{r=1}^j \{X_{r:k} \geq \mu_0 - d_{j:k}^{(1)} \frac{\sigma}{\sqrt{n}}\}\right) \\ &\leq P_\mu\left(\bigcup_{j=1}^{k-i+1} \bigcup_{r=1}^j \{Z_{r:k} + \frac{\sigma r \tau_0}{\sqrt{n}} \geq -d_{j:k}^{(1)}\}\right) \end{aligned}$$

which is decreasing in μ_r , $r = 1, \dots, k-i+1$.

Hence

$$\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i+1}) \geq P(Z_{k-i+1:k} \geq -d_{k-i+1:k}^{(1)})$$

On the other hand,

$$\begin{aligned} & \inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i+1}) \\ &= P_\mu^*(\bigcup_{j=1}^{k-i+1} \{X_{j:k} \geq \mu_0 - d_{j:k}^{(1)} \frac{\sigma}{\sqrt{n}}\}) \\ &= P(Z_{k-i+1:k} \geq -d_{k-i+1:k}^{(1)}) \end{aligned}$$

whenever $\mu^* = (\underbrace{-\tau_0, -\tau_0, \dots, -\tau_0}_{i}, \underbrace{\tau_0, \dots, \tau_0}_{k-i}) \in \mathcal{C}_i$

Thus, we have

$$\inf_{\mu \in \mathcal{C}_i} P_\mu(CS|_{i+1}) = P(Z_{k-i+1:k} \geq -d_{k-i+1:k}^{(1)}).$$

Since $\hat{z}_{k-i+1:k} = \min\{z_{k-i+1}, \dots, z_{k-i+1+i}\}$ have the same distributions as

$$\hat{z}_{1:i} = \min\{z_1, \dots, z_i\},$$

let

$$\begin{aligned} V_i &= \hat{z}_{1:i} \\ &= \min_{1 \leq j \leq i} \frac{1}{r} \sum_{j=1}^r z_j, \end{aligned} \quad (1.3.4)$$

we have

$$\inf_{x \in \mathbb{R}} P_p(CS|x_1) = P(V_i \sim -d_{k-i+1:k}^{(1)}), \quad i = 1, \dots, k. \quad (1.3.4)$$

Theorem 1.3.1. In case I, (ω_0 known, common known σ^2 and common sample size n), if $d_{k-i+1:k}^{(1)}$ is the solution of equation

$$P(V_i \sim -x) = p^*, \quad (1.3.5)$$

where

$$V_i = \min_{1 \leq j \leq i} \frac{1}{r} \sum_{j=1}^r z_j \text{ and } z_i \text{ are i.i.d. } F(\cdot),$$

$i = 1, \dots, k$ then x_1 satisfies the P^* -condition.

Proof. For any i , $1 \leq i \leq k$,

$$\inf_{x \in \mathbb{R}} P_p(CS|x_1) = P(V_i \sim -d_{k-i+1:k}^{(1)}) = p^*,$$

so x_1 satisfies the P^* -condition.

Therefore, the problem of finding the $d_{i:k}^{(1)}$'s reduced to finding the distributions of V_1, \dots, V_k . This is achieved by using some theorems in the theory of random walk.

1.3.3. Some Theorems in the Theory of Random Walk

Suppose Y_1, Y_2, \dots are independent random variables with a common distribution B not concentrated on a half-axis, i.e. $0 < P(Y_1 = 0) < 1$, $P(Y_1 > 0) > 1$. The induced random walk is the sequence of random variables

$$S_0 = 0, S_n = Y_1 + \dots + Y_n, \quad n = 1, 2, \dots$$

Let

$$\tau_n = P(S_1 \geq 0, \dots, S_{n-1} \geq 0, S_n \geq 0) \quad (1.3.6)$$

and

$$\varphi(s) = \prod_{n=1}^{\infty} \tau_n s^n, \quad 0 < s \leq 1. \quad (1.3.7)$$

Then we have the following theorem which was discovered by Anderson (1953). Feller (1971) gave an elegant short proof.

Theorem 1.3.2.

$$\log \frac{1}{\tau_n}(s) = \sum_{n=1}^{\infty} \frac{s^n}{n} P(S_n = 0). \quad (1.3.8)$$

Theorem 1.3.3. (Feller (1971))

Let

$$p_n = P(S_1 = 0, \dots, S_n = 0),$$

then

$$p(s) = \prod_{n=1}^{\infty} p_n s^n = \frac{1}{\tau_n}(s), \quad (1.3.9)$$

hence

$$\log p(s) = \sum_{n=1}^{\infty} \frac{s^n}{n} p(S_n = 0). \quad (1.3.10)$$

By symmetry, the probabilities

$$q_n = P(S_1 \leq 0, \dots, S_n \leq 0) \quad (1.3.11)$$

have the generating function q given by

$$\log q(s) = \sum_{n=1}^{\infty} \frac{s^n}{n} P(S_n \leq 0). \quad (1.3.12)$$

Note: The above two theorems remain valid if the signs $-$ and $+$ are replaced by \sim and \ll , respectively.

Now, let

$$U_j^r = \max_{1 \leq i \leq j} \frac{1}{r} \sum_{i=1}^r Z_i^r, \quad j = 1, 2, \dots, \quad (1.3.13)$$

and

$$V_j^r = \min_{1 \leq i \leq j} \frac{1}{r} \sum_{i=1}^r Z_i^r, \quad j = 1, 2, \dots, \quad (1.3.14)$$

where Z_i^r 's are i.i.d. with absolutely continuous c.d.f. $G(\cdot)$. We would like to apply Theorem 1.3.3 to get the distribution of U_j^r and V_j^r , $j = 1, 2, \dots$.

Remark 1.3.1. The distribution of U_j^r , $j = 1, \dots, k$ for some $k < r$, will be used whenever our goal is changed to select a subset containing no population with parameter smaller than the control.

Theorem 1.3.4. The generating function $q(s)$ of $P(U_j^r \leq x)$, $j = 1, \dots$ is

$$\prod_{j=1}^{r-1} P(U_j^r \leq x) = e^{-x} \prod_{n=1}^{\infty} \frac{1}{n!} s^n P(Z_n \leq 0) = (1 - F_x)^s$$

where

$$S_n = \sum_{i=1}^n (Z'_i - x), \quad n = 1, 2, \dots,$$

if the distribution of $Y_1 = Z'_1 - x$ is not concentrated on a half-axis.

Proof. Since the distribution of random variable $Y_i = Z'_i - x$ is not

concentrated on a half-axis, and Y_i 's are i.i.d. let $S_r = \sum_{i=1}^r (Z'_i - x)$,
 $r = 1, \dots, k$. Then

$$(U_j > x) \geq \max_{1 \leq r \leq j} \frac{1}{r} S_r \geq 0 \Leftrightarrow (S_1 \geq 0, \dots, S_j \geq 0).$$

By Feller's Theorem 1.3.3, we complete the proof.

Similarly, $(V_j > x) = (S_i \geq 0, \quad i = 1, 2, \dots, j)$,

where

$$S_i = \sum_{r=1}^i (Z'_r - x).$$

Theorem 1.3.5. The generating function $p(s)$ of $P(V_j > x)$ is

$$\sum_{j=1}^{\infty} s^j P(V_j > x) = \exp \left[\sum_{n=1}^{\infty} \frac{1}{n} s^n P(S_n > 0) \right], \quad (1.3.16)$$

if the distribution of $Y_1 = Z'_1 - x$ is not concentrated on a half-axis.

Corollary 1.3.1. Both Theorem 1.3.4 and Theorem 1.3.5 hold for all
 x such that $0 < G(x) < 1$.

Proof. Let $Y_1 = Z_1^t - x$, then

$$P(Y_1 < 0) = G(x)$$

and

$$0 < G(x) < 1,$$

hence Y_1 is not concentrated on a half-axis.

Corollary 1.3.2. Both Theorem 1.3.4 and Theorem 1.3.5 hold for all x whenever $G = \phi$, c.d.f. of $N(0,1)$, or $G = F$ which is defined at the beginning of Section 1.3.

Proof. Followed immediately by Corollary 1.3.1.

Note that in the case of location parameter of normal population,

$$P(U_n^t < -x) = P(V_n^t > x).$$

Let

$$\gamma_j(x) - \gamma_j = P(S_j < 0), \quad j = 1, 2, \dots,$$

$$a(s) = \frac{1}{n} \sum_{j=1}^n \gamma_j^n,$$

we have

$$p(s) = \prod_{j=1}^n s^{j-1} P(V_j^t < x) = \exp(a(s)).$$

$$\text{Lemma 1.3.1. } p^{(n+1)}(s) = \sum_{j=0}^n \binom{n}{j} p^{(j)}(s) a^{(n+1-j)}(s), \quad n.$$

Proof. Since $p'(s) = p(s) + a'(s)$, the result can be proved by induction on n .

Theorem 1.3.6. Under the assumption of Theorem 1.3.5

$$\begin{aligned} P(V_{n+1}' \leq x) &= \frac{1}{(n+1)!} \lim_{s \rightarrow 0^+} \frac{d^{n+1}}{ds^{n+1}} p(s) \\ &= \frac{1}{n+1} \sum_{j=0}^n P(V_j' \leq x) \gamma_{n-j+1}, \quad n = 0, 1, 2, \dots \quad (1.3.17) \end{aligned}$$

where

$$P(V_0' \leq x) = 1, \quad \forall x.$$

Proof. By Lemma 1.3.1, we have

$$\begin{aligned} P(V_{n+1}' \leq x) &= \frac{1}{(n+1)!} \lim_{s \rightarrow 0^+} p^{(n+1)}(s) \\ &= \frac{1}{n+1} \sum_{j=0}^n \frac{1}{(n+1)!} \frac{n!}{j!(n-j)!} p^{(j)}(0) [(n-j)!]^{-1} \gamma_{n+1-j} \\ &= \frac{1}{n+1} \sum_{j=0}^n \frac{p^{(j)}(0)}{j!} \gamma_{n+1-j} \\ &= \frac{1}{n+1} \sum_{j=0}^n P(V_j' \leq x) \gamma_{n+1-j}. \end{aligned}$$

Similarly, we have

$$P(U_{n+1}' \leq x) = \frac{1}{n+1} \sum_{i=0}^n P(U_{n-i+1}' \leq x) P(S_i = 0). \quad (1.3.18)$$

1.3.4. Limiting Distributions of U_n' and V_n'

Let $f_n(x) = P(U_n' \leq x)$ and $F_n(x)$ denote the limiting distribution function as $n \rightarrow \infty$ of U_n' . Suppose the distribution of random variable $Y_1 - Y_1' + x$ is not concentrated on a half axis, then we have

$$1 - F_n(x) = P(S_1 > 0) + \sum_{r=2}^n P(S_1 > 0, \dots, S_{r-1} = 0, S_r = n),$$

$$1 - F_n(x) = \lim_{s \downarrow 1^-} \varphi(s),$$

and apply Andersen-Feller Theorem 1.3.2, we have

$$F_n(x) \approx \exp \left(- \sum_{r=1}^{k-1} \frac{1}{r} P(S_r > 0) \right). \quad (1.3.1)$$

Similarly,

$$G_n(x) \approx \exp \left(- \sum_{r=1}^{k-1} \frac{1}{r} P(S_r < 0) \right). \quad (1.3.2)$$

where

$$G_n(x) = P(V_n < x).$$

Let

$$G_n(-d_{1:n}^{(1)}) = P^*. \quad (1.3.3)$$

If Z_i , $i = 1, \dots, k$, are independent identically distributed $N(0,1)$, then we can use the recurrence formula of Theorem 1.1.6 to solve the equations $P(V_i < -d_{k-i+1:k}) = P^*$, $i = 1, \dots, k$. Hence in Case I, $\varphi_j(x) = (-x)^j$.

Remark 1.3.2. From formula (1.3.3) we know that $d_{k-i+1:k}^{(1)}$ ($i = 1, \dots, k$) does not depend on k . And we have $d_{k-i+1:k}^{(1)} = d_{1:i}^{(1)}$. These values for $k = 1$ (1) 6, 10, ... and $P^* = .99, .975, .95, .925, .9, .85, .75, .7, .65$ are tabulated in Table I.

1.3.5. Some Other Forms of Selection Procedure 1

Lemma 1.3.2. $d_{1:i}^{(1)}$ is increasing in i .

Proof. By Remark 1.3.2 and the fact

$$v_{i+1} = \min(v_i, \frac{iv_i + z_{i+1}}{i+1}).$$

Lemma 1.3.3. If c_j , $1 \leq j \leq i \leq k$ is decreasing in j , then

$$\frac{\bigcup_{j=1}^i x_{j:k}}{i} - c_j = \frac{\bigcup_{j=1}^{i-1} x_{j:k}}{i-1} - c_j$$

$$\text{Proof. } \frac{\bigcup_{j=1}^i x_{j:k}}{i} - c_j = \frac{\bigcup_{j=1}^{i-1} x_{j:k}}{i-1} - c_j,$$

since

$$x_{j:k} < \hat{x}_{j:k}, \quad 1 \leq j \leq k.$$

On the other hand, if

$$x_{r:k} = c_r \quad \text{for some } r, \quad 1 \leq r \leq i$$

then

$$x_{s:k} = c_r \quad \text{for some } s, \quad 1 \leq s \leq r,$$

since

$$x_{r:k} = \max_{1 \leq s \leq r} x_{s:k}.$$

Because c_j is decreasing in j , this implies $x_{s:k} = c_s$ for some s , $1 \leq s \leq r$.

Hence we have

$$\sum_{j=1}^i (x_{j:k} - c_j) = \sum_{j=1}^i (x_{j:k} - c_j'),$$

therefore the lemma is proved.

Definition 1.3.2. We define a selection procedure \cdot_1' by replacing the inequality in the i th step of procedure \cdot_1 by the inequality

$$x_{i:k} \geq n - d_{i:k}' \quad i = 1, \dots, k$$

where $d_{i:k}', \dots, d_{k:k}'$ are the smallest values such that \cdot_1' satisfies the P^* -condition.

Theorem 1.3.7. The selection procedure \cdot_1 and \cdot_1' are identical since $d_{i:k}^{(1)} = d_{i:k}', i = 1, \dots, k$.

Proof. For any i , $1 \leq i \leq k$, by Theorem 1.3.1

$$P^* = \inf_{\cdot \in \cdot_1} P_{\cdot}(CS^{(1)}_1) = P(Z_{k-i+1:k} \geq d_{k-i+1:k}^{(1)}).$$

On the other hand, using the same arguments as Section 1.3.7, we have

$$P^* = \inf_{\cdot \in \cdot_1'} P_{\cdot}(CS^{(1)}_1) = P(Z_{k-i+1:k} \geq d_{k-i+1:k}').$$

Hence we have $d_{i:k}^{(1)} = d_{i:k}', i = 1, \dots, k$.

Since $x_{1:k} = x_{i:k}$, the first step of \cdot_1 and \cdot_1' are identical. For $i = 2, \dots, k$, the event

$$\begin{aligned}
 & \text{select } i_1, \dots, i_{k-1} \leftarrow \inf_{\substack{0 \leq x_{j;k} \leq d \\ j=1}} \{x_{j;k} : 0 \leq x_{j;k} \leq d\}_{j=1}^{k-1} \\
 & \quad \vdots \\
 & \quad \text{select } i_1, \dots, i_{k-1} \leftarrow \inf_{\substack{0 \leq x_{j;k} \leq d \\ j=1}} \{x_{j;k} : 0 \leq x_{j;k} \leq d\}_{j=1}^{k-1} \\
 & \quad \vdots \\
 & \quad \text{select } i_1, \dots, i_{k-1}
 \end{aligned}$$

by Lemma 1.3.2 and Lemma 1.3.3. Hence selection procedures τ_1 and τ_2 are identical.

1.3.6. Some Other Proposed Selection Procedures τ_2, τ_3, τ_4

In Case I, we proposed some other selection procedures:

Definition 1.3.3. We define a selection procedure τ_3 by

$$\tau_3: \text{select } i_1 \text{ if and only if } x_{i;k} \geq 0 \geq d_{i;n}, \quad i = 1, \dots, k$$

where d is the smallest value such that τ_3 satisfies the P*-condition.

Theorem 1.3.8. Under assumptions of Case I, and selection procedure τ_3 , if we select population i_1 , then we will select populations i_j ,

for all $j = 1$.

Proof. Since $x_{i;k} \geq x_{j;k}$ for all $j = 1$.

Evaluation the Value of τ_3

For any $i_1, 1 \leq i_1 \leq k$, we have

$$\begin{aligned}
 \inf_{\substack{0 \leq x_{j;k} \leq d \\ j \neq i_1}} P(\tau_3) & \leq \inf_{\substack{0 \leq x_{j;k} \leq d \\ j \neq i_1}} P(\tau_{k-i+1;k} \geq 0 \geq d)_{j=1}^{k-1} \\
 & \leq P(V_{i_1} \geq -d)
 \end{aligned}$$

by the same argument for selection procedure π_1 and here

$$Y_i = Z_{1:i} - \min_{1 \leq j \leq i} \frac{1}{r} \sum_{j=1}^r Z_{j:i}$$

We need the constant d such that $P(Y_i \geq -d) \geq P^*$ holds for all i ,

$1 \leq i \leq k$. By Lemma 1.3.2 we have $d = d_{1:k}^{(1)}$. Hence we have the following theorem.

Theorem 1.3.3. Selection procedure π_p satisfies the P^* -condition with $d = d_{1:k}^{(1)}$.

Corollary 1.3.3. If S_1 and S_2 are the selected subsets associated with selection procedures π_1 and π_2 , respectively, then $\pi_1 = \pi_2$.

Proof. Proof follows from Lemma 1.3.2.

Definition 1.3.4. The procedure π_3 is defined as follows:

Step 1. Select π_1 , $i = 1$ and stop, if

$$X_1 - \mu_0 - d_1 \frac{\sigma}{\sqrt{n}} > 0,$$

otherwise reject π_1 and go to step 2.

Step 2. Select π_2 , $i = 2$ and stop, if

$$X_2 - \mu_0 - d_2 \frac{\sigma}{\sqrt{n}} > 0,$$

otherwise reject π_2 and go to step 3.

⋮

Step $k-1$. Select π_k , $i = k-1$ and stop, if

$$X_{k-1} - \mu_0 - d_{k-1} \frac{\sigma}{\sqrt{n}} > 0,$$

otherwise reject x_{k+1} and go to step k.

Step k. Select x_k and stop, if

$$x_k - \bar{x}_0 \geq d_k \frac{\sigma}{\sqrt{n}},$$

otherwise reject x_k .

Here $\bar{x}_j = \max(x_1, \dots, x_j)$ and d_j 's are the smallest values such that x_j satisfies the P*-condition.

Evaluation of d_j 's

For any i , $1 \leq i \leq k$,

$$\begin{aligned} \inf_{u \in C_i} P_u(CS^{1,3}) &= \inf_{u \in C_i} P_u(\bar{x}_{j+1} - x_i - u_0 \geq d_j \frac{\sigma}{\sqrt{n}}) \\ &= P_u(Z_{k-i+1} \geq d_{k-i+1} \frac{\sigma}{\sqrt{n}}) \\ &= P(Z_{k-i+1} \geq d_{k-i+1}) \\ &= P^*, \quad Z_i \sim F(\cdot; 0, 1). \end{aligned}$$

This implies $d_{k-i+1} = d$ for all i , and

$$\begin{aligned} d &= F^{-1}(1-P^*), \\ &= F^{-1}(P^*), \quad \text{if } F \text{ is symmetric} \\ &= F^{-1}(P^*), \quad \text{if } X_i \sim N(\mu_i, \sigma^2/n). \end{aligned}$$

Similar to the selection procedure γ_1 , we have the following theorem:

Theorem 1.3.10. Selection procedure γ_3 satisfies the P*-condition with

$$d_j = F^{-1}(1-P^*).$$

Definition 1.3.5. Selection procedure γ_1 is defined as follows:

Step 1. Select x_1 , $i = 1$ and stop, if

$$x_1 - \mu > d_{\frac{\alpha}{\sqrt{n}}},$$

otherwise reject x_1 and go to step 2.

Step 2. Select x_2 , $i = 2$ and stop, if

$$x_2 - \mu > d_{\frac{\alpha}{\sqrt{n}}},$$

otherwise reject x_2 and go to step 3.

⋮

⋮

Step k-1. Select x_i , $i = k-1$ and stop, if

$$x_{k-1} - \mu > d_{\frac{\alpha}{\sqrt{n}}},$$

otherwise reject x_k .

Here

$$\begin{aligned}d &= F^{-1}(1-P^*) \\&= F^{-1}(P^*) \text{ if } F \text{ is symmetric.}\end{aligned}$$

Theorem 1.3.11. The selection procedure γ_1 satisfies the P* condition.

Proof. For any i , $1 \leq i \leq k$,

$$\inf_{x_i} P_{\mu}(CS^{(i)}) = P_{\mu}(x_{k+1}) = 1 - P^*.$$

Theorem 1.3.12. The selection procedures γ_1 and γ_2 are identical.

Proof. The proof is simple hence it is omitted.

The following procedure π_4 was given by Gupta and Sobel (1954), without assuming any ordering prior:

Definition 1.3.6. The selection procedure π_4 is defined as follows:

π_4 : Select x_i if and only if $x_i - d \geq 0$, $i = 1, \dots, k$

where d is the smallest constant such that π_4 satisfies the P*-condition.

It was shown that the value d is determined by the equation

$$F(-d) = \frac{1}{1 + P^k}$$

or

$$F(d) = P^k \text{ if } F \text{ is symmetric.}$$

1.3.7. A Dual Problem

We start with the same assumptions as in Section 1.3.1 Case 1, but change our goal to select a subset which contains no bad populations; the definition of a correct selection (CS) will now be changed to select a subset that contains no bad populations and the P*-condition will be defined based on this new definition of correct selection (CS).

In location parameter case, this problem is a dual problem of the original problem, namely, "select a subset which contains all good populations under ordering prior assumption".

One method to solve this problem is that, first, change the signs of all statistics and the control to opposite signs; then use a procedure for selecting a subset which contains all "new good" populations.

where the "new good" populations are the "old bad" populations before changing signs; finally, reject the selected subset and keep the remainders as the desired selected subset. Let ω_i , $i = 1, 2, \dots, 4$ denote the above procedure which corresponds to ω_i , $i = 1, 2, \dots, 4$, respectively.

Theorem 1.3.3. The selection procedure ω_i , $i = 1, 2, \dots, 4$ satisfies the P^* -condition in which the correct selection (CS) means that it selects a subset which contains no bad population.

Proof. Given P^* and observations, for any selection procedure ω_i , $i = 1, 2, \dots, 4$, after changing the signs of all associated states, the probability that the selected subset S contains all "new good" populations is not less than P^* . If we reject the selected subset S , then the complement subset S^C of S contains any "new good" populations with probability less than $1-P^*$, but the "new good" populations are the originally bad populations so what we have is that the subset S^C contains any originally bad population with probability less than $1-P^*$, in other words, subset S^C contains no bad population with probability greater than or equal to P^* . Since this is true for all arbitrary true configurations, we have completed the proof.

Remark 1.3.3. It is easy to see that the value $d_{i+1}^{(k)} e^{-\lambda t}$ of ω_1 which was used by ω_1 in the i th step is determined by the equation

$$P(\omega_{k-i+1} + d_{i+1}^{(k)} e^{-\lambda t}) = P^*,$$

where

$$\frac{P(\omega_{k-i+1} + d_{i+1}^{(k)} e^{-\lambda t})}{P(\omega_{k-i+1})} = \frac{e^{-\lambda t}}{P(\omega_{k-i+1}) - P(\omega_{k-i+1} + d_{i+1}^{(k)} e^{-\lambda t})}.$$

If the distribution F is symmetric, then

$$d_{i;k}(.)_1 = d_{i;k}^{(1)}.$$

1.3.8. Some Proposed Selection Procedures $\psi_i^{(2)}$, $i = 1, 2, 3, 4$

When θ_0 is Unknown

Case II. θ_0 unknown, common σ^2 known, common sample size n .

Definition 1.3.7. We define a selection procedure $\psi_1^{(2)}$ by requiring the inequalities

$$x_{i;k} - x_0 \geq d_{i;k}^{(1)} \frac{1}{n}, \quad i = 1, \dots, k$$

in procedure ψ_1 (Definition 1.3.1) with

$$x_{i;k} - x_0 \geq d_{i;k}^{(2)} \frac{1}{n}, \quad i = 1, \dots, k, \text{ respectively.}$$

Here $x_0 = \frac{n}{i+1} x_{0i}/n$, $d_{i;k}^{(2)}$, $i = 1, \dots, k$ are the smallest constants such

that the selection procedure $\psi_1^{(2)}$ satisfies the P*-condition.

Similar to the Case I, we have the following theorem:

Theorem 1.3.14. For any $t_1 < t_2 < \dots < t_k < t_{k+1}$, is determined by the equation

$$\int_{t_1}^{t_2} F(V_1 - t) + d_{k-i+1;k}^{(2)} dF(t) = P^*, \quad (1.3.14)$$

It is easy to see that $d_{k-i+1:k}^{(2)} > d_{1:i}^{(2)}$. The following theorem gives us an identical form of the selection procedure $\psi_1^{(2)}$.

Theorem 1.3.15. The selection procedure $\psi_1^{(2)}$ will not be characterized if the statistics $X_{i:k}$, $i = 1, \dots, k$, are replaced by $\bar{X}_{i:k}$, $i = 1, \dots, k$, respectively.

Proof. The proof is the same as that in Case I and hence it is omitted.

The values $d_{1:i}^{(2)}$, $i = 1, \dots, k$ are tabulated in Table II for $k = 1 (1) 6, 8, 10$, and $P^* = .99, .975, .95, .925, .90, .85, .75, .70, .65$.

Similar to the Case I, we propose a selection procedure $\psi_2^{(2)}$ as follows:

Definition 1.3.8. We define a selection procedure $\psi_2^{(2)}$ by

$\psi_2^{(2)}$: Select τ_i if and only if $\bar{X}_{i:k} - \bar{X}_0 = d_{1:n}^{(2)} + 1 - E_{i:k}$

where d is the smallest value such that $\psi_2^{(2)}$ satisfies the P^* -condition. Then, similar to Theorem 1.3.9 we have:

Theorem 1.3.16. Under assumptions of Case II, the selection procedure $\psi_2^{(2)}$ satisfies the P^* -condition with $d = d_{1:n}^{(2)}$.

Next, we define a selection procedure $\psi_3^{(2)}$ which is identical to $\psi_2^{(2)}$ but replace \bar{X}_0 by $\bar{\tau}_0$, the sample mean of population τ_0 .

Definition 1.3.9. The selection procedure $\{x_i^{(2)}\}$ is defined by condition
 $x_i = x_0 + d_i \frac{x_i - x_0}{\sqrt{n}}$ in γ_3 (Definition 1.3.4) by $x_i = x_0 + d_{i+1}^* \frac{x_i - x_0}{\sqrt{n_i}}$, $i = 1, \dots, k$
where d_1^*, \dots, d_k^* are the smallest values such that $\{x_i^{(2)}\}$ satisfies the
P*-condition.

Similar to Theorem 1.3.10 we have:

Theorem 1.3.12. The selection procedure $\{x_i^{(2)}\}$ satisfying the P*-condition
with $d_i = d$, $i = 1, \dots, k$ where d is determined by the equation

$$\int_{-\infty}^d [1 - F(t-d)]dF(t) = P^*, \quad (1.3.12)$$

$$\int_{-\infty}^d F(d-t)dF(t) = P^*, \text{ if } F \text{ is symmetric.}$$

And $\{x_i^{(2)}\}$ will not be changed if the statistics x_i is replaced by \bar{x}_i ,
the sample mean of population γ_i for $i = 1, \dots, k$.

The following selection procedure $\{x_i^{(3)}\}$ was proposed by Guntar and
Sobel (1958):

Definition 1.3.10. The selection procedure $\{x_i^{(3)}\}$ is defined by

$\{x_i^{(3)}\}$: Select $x_i^{(3)}$ if and only if $x_i = x_0 + d \frac{x_i - x_0}{\sqrt{n_i}}$, $i = 1, \dots, k$

where d is determined by the following equation if F is normal distri-
bution:

$$\sum_{i=1}^k \left[F\left(\frac{x_i - x_0}{\sqrt{n_i}} + d\right) \right] F(u)du = P^*, \quad (1.3.13)$$

For the special case $n_j = n$ ($j = 0, 1, \dots, k$)

$$\int_{-\infty}^{\infty} F^k(t+d)f(t)dt = P^*. \quad (1.3.25)$$

If F is normal distribution $N(0,1)$, the tables of d -values satisfying the Equation (1.3.25) for several values of P^* are given in Bechhofer (1954) for $k = 1, 2, 3, 4$ and in Gupta (1946) for $k = 1$.

1.3.9. Some Proposed Selection Procedures $\{S_i^{(j)}\}$, $i = 1, \dots, n$

When common Variance σ^2 is unknown

Case III. μ_0 known, common variance σ^2 unknown, $n_1 = n - 1$

In this case, we assume that $(X_i - \mu_0)^2$ which is the square of $N(0,1)$.

Definition 1.3.11. We define the selection procedure $\{S_i^{(j)}\}$ to replace the inequalities

$$c_{ijk} = n + d_{ijk} \frac{1}{n-1}, \quad i = 1, \dots, k$$

in procedure $\{S_i\}$ (Definition 1.3.10) by

$$S_{ijk}^{(j)} < S_{ijk+1}^{(j)}, \quad i = 1, \dots, k, \text{ respectively,}$$

where $d^{(j)}$'s are the smallest values such that $S_{ijk}^{(j)}$ satisfies the P^* -condition, $S_i^{(j)}$ denotes the i -th observation of X_i for $i = 1, \dots, k(n-1)$, that is,

$$S_{ijk}^{(j)} = \frac{X_{ijk} - \mu_0}{\sqrt{\frac{1}{n-1} \sum_{i=1}^{n-1} (X_{ijk} - \mu_0)^2}}.$$

Note that $\frac{\chi^2_{N-k}}{k}$ has the chi-square distribution χ^2 with k degrees of freedom.

By using similar arguments as in Case I, we have:

Theorem 1.3.18. The equation which determines the constant $d_{k-i+1;k}^{(1)}$ is

$$P(V_i \leq d_{k-i+1;k}) = P^* \quad (1.3.18)$$

or

$$\int_0^{V_i} P(V_i \leq d_{k-i+1;k}) q_i(y) dy = P^* \quad (1.3.18)$$

where

$$V_i = \min_{1 \leq r \leq i} \frac{1}{r} \sum_{j=1}^r z_j$$

and $q_i(y)$ is the density of $\frac{S_i - \bar{z}_i}{\sqrt{t}}$.

We can rewrite formula (1.3.18) as

$$\int_0^{V_i} P(V_i \leq d_{k-i+1;k}) \sqrt{t} dt e^{-\frac{V_i^2}{t}} = P^* \quad (1.3.18)$$

or

$$\int_0^{V_i} P(V_i \leq d_{k-i+1;k}) \sqrt{t} dt e^{-\frac{V_i^2}{t}} = \frac{1}{2} \int_0^{V_i^2} e^{-\frac{t}{2}} dt = P^* \quad (1.3.18)$$

Remark 1.3.4. The values of $d_{k-i+1;k}^{(3)}$, $i = 1, \dots, k$ depend on $k+i-1$, hence $d_{k-i+1;k}^{(3)} \neq d_{1;i}^{(3)}$.

By using Rabinowitz and Weiss table (1959) (with $n = 3, 4, \dots, 10$), we have evaluated and tabulated the values of $A_{k-i+1}^{(2)}$, $i = 1, \dots, k$, in Table III, for $k = 2, 4, 6, 8$, $\alpha^* = .29, .495, .745, .945, .995, .998, .999$, .80, and .75 with common sample size $n = 3, 5, 9$, and 13.

For $k = 6$ and $n = 24$, we can take $\{ \frac{1}{2} \}_{k=1}^6$ as an appropriate choice of $\delta_{k+1, k+2}^{(3)}$.

Definition 1.3.11. We define the selection procedure $\hat{f}_n^{(k)}$ by

(3): Select j if and only if $x_{ijk} - d^{(j)} \leq s$ for all i, k , where s is defined as in procedure (3), and $d^{(j)}$ is the smallest constant such that (3) satisfies the P*-condition.

As before, it can be shown that $\frac{d\psi}{dx} = \frac{d\psi}{dp}$.

Remark 1, 2, 5. Theorem 1, (1,2) still holds, for case III, i.e. the selection procedure $\hat{\gamma}_j^{(t)}$ will not be changed if we replace the F-test statistics τ_{ijk} by ζ_{ijk} , respectively. But this is not true in general, true for selection procedure $\hat{\gamma}_j^{(t)}$.

Definition 1.3.1. The selection procedure $\tilde{S}_n^{(k)}$ is defined to have the same form as procedure $S_n^{(k)}$, except that the inequality defining the i th step of procedure $\tilde{S}_n^{(k)}$ is replaced by

$$f_1 = \frac{1}{2} \left(-\partial_x^2 + \partial_y^2 + \partial_z^2 + \partial_w^2 \right) f_0 + \text{higher order terms}$$

The proof of the following theorem uses the same arguments as that in Case I, hence it is omitted.

Theorem 1.3.19. The equation which determines the constant d of selection procedure $\tau_3^{(3)}$ is

$$\int_0^{\infty} \tau_3(yd) q_{\tau_3}(y) dy = P^*, \quad (1.3.31)$$

Gupta and Sobel (1958) gave a selection procedure $\tau_4^{(4)}$ in this case. It is as follows:

$\tau_4^{(4)}$: Select x_i if and only if $x_i > c_0 + d \sum_{j=1}^k \frac{S_j}{m_j}$, $i = 1, \dots, k$

and the equation which determines d is

$$\int_0^{\infty} \tau_4^k(yd) q_{\tau_4}(y) dy = P^*. \quad (1.3.32)$$

1.3.10. Some Proposed Selection Procedures $\tau_i^{(4)}$, $i = 1, 2, 3, 4$

When Both Control c_0 and Common Variance σ^2 are Unknown.

Case IV. c_0 unknown, common variance σ^2 unknown and common sample size n .

We assume that in this case distribution F is the c.d.f. $N(t, 1)$, and denoted by τ . We replace c_0 in each selection procedure $\tau_i^{(3)}$ by $c_0 + t$, $t \geq 0$, and get four procedures $\tau_i^{(4)}$, $1 \leq i \leq 4$, respectively. Let $\chi^2(t)$ denote the c.d.f. of the chi-square distribution with $k(n-1)$ degrees of freedom. The constant $d_{k-i+1;k}^{(4)}$, $i = 1, \dots, 4$, of procedure $\tau_1^{(4)}$ is determined by

$$\int_0^{\infty} \int_{-\infty}^{\infty} P(Y_i = u + d_{k-i+1, k-i}^{(4)} t) d(u) dt, \quad i = 1, \dots, k.$$

The constant d of procedure $\psi_2^{(4)}$ is

$$d = d_{1, k}^{(4)},$$

The constants d of procedures $\psi_3^{(4)}$ and $\psi_4^{(4)}$ are determined by

$$\int_0^{\infty} \int_{-\infty}^{\infty} r(u + td) c'(u) d(u)^2(t) = p, \quad i = 3, 4,$$

with $r = 1$ and k , respectively, and their values for selected values of P^* , k and i are given in Gupta and Sebel (1967) and Bennett (1971).

1.3.11. Properties of the Selection Procedures

Under simple ordering prior, it is natural to require that an order selection procedure is order-preserving as defined below:

Definition 1.3.14. A selection procedure ψ is order-preserving if it selects γ_j with parameter α_j , and if $\gamma_j < \gamma_s$, then it also selects γ_s . Procedure ψ is weak order-preserving or monotone if

$$P(\gamma_j \text{ is selected}) \geq P(\gamma_j \text{ is selected}) \text{ whenever } \gamma_j < \gamma_s.$$

It is easy to see that any order-preserving selection procedure is weak order-preserving, but the converse is not true.

Now, let $\psi_i^{(1)} = \gamma_i$, $i = 1, \dots, k$.

Theorem 1.3.20. The selection procedures $\psi_1^{(1)}$, $\psi_2^{(1)}$ and $\psi_3^{(1)}$ are

order-preserving and procedure $\psi_4^{(1)}$ is monotone for $i = 1, \dots, k$.

Proof. The proof follows immediately from the definition of the procedures.

Given observations $X = x = (x_0, \dots, x_k)$ where x_i is the sample mean of population γ_i , $i = 1, \dots, k$, and $x_0 = \bar{x}_0$ if \bar{x}_0 is known, otherwise x_0 is the sample mean of population γ_0 . Let

$$\gamma_i(x, \cdot) = P(\gamma_i \text{ included in the selected subset } x = x_0, \dots, x_k) \\ \text{for } i = 1, \dots, k.$$

Definition 1.3.16. A selection procedure γ is called translation-invariant if for any $c \in \mathbb{R}^{k+1}$, $c \in \mathbb{R}$

$$\gamma(x_0 + c, x_1 + c, \dots, x_k + c, \cdot) = \gamma(x_0, \dots, x_k; x_0 + c, \dots, x_k + c)$$

Theorem 1.4.21. The selection procedures $\gamma_1^{(i)}, \gamma_2^{(i)}, \gamma_3^{(i)}$ and $\gamma_4^{(i)}$ are translation-invariant for $i = 1, 2, 3, 4$.

Proof. By Corollary 1.3.1 the isotonic regression is a linear operator. On the other hand,

$$\frac{\sum_{i=1}^n \gamma_i(x_i + c)}{n} = \frac{\sum_{i=1}^n \gamma_i(x_i)}{n} + c$$

hence we have the result.

Expected Number (Size) of Bad Populations in the Selected Subset

Suppose the control γ_0 is known and we have common sample size n and common known variance σ^2 without loss of generality we assume that $\gamma_0 = 0$ and $\gamma_0 = 1$. Let $\gamma(x)$ denote the expected number of bad populations in the selected subset x using the selection procedure

, then for any i , $0 \leq i \leq k$,

$$\begin{aligned} & \sup_{\mu \in \mathcal{K}_{k-i}} E(S^{(1)}_{\mu, 1}) \\ &= \sup_{\mu \in \mathcal{K}_{k-i}} \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\mu, j} \cap d_{1;k}^{(1)}) \\ &= \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\mu, j} \cap d_{1;k}^{(1)}) \quad (\text{from (1.3.35)}) \end{aligned}$$

On the other hand, for procedure φ

$$\sup_{\mu \in \mathcal{K}_{k-i}} E(S^{(1)}_{\mu, 1}) = \sum_{r=1}^i P(\bigcup_{j=1}^r Z_{\varphi, j} \cap d_{1;k}^{(1)}) \quad (\text{from (1.3.35)})$$

Formula (1.3.36) is increasing in i and is greater than or equal to

Formula (1.3.35), since

$$d_{1;k}^{(1)} \geq d_{1;k+1}^{(1)} \geq d_{1;k}^{(1)}$$

Therefore, we have the following theorem.

Theorem 1.3.22. For any i , $0 \leq i \leq k$

$$\sup_{\mu \in \mathcal{K}_i} E(S^{(1)}_{\mu, 2}) \geq \sup_{\mu \in \mathcal{K}_i} E(S^{(1)}_{\mu, 1}),$$

$$\sup_{\mu \in \mathcal{K}_0} E(S^{(1)}_{\mu, 2}) \geq \sup_{\mu \in \mathcal{K}_0} E(S^{(1)}_{\mu, 1}).$$

Theorem 1.3.23. In Section 1.3.1, Case 1, for any i , $0 \leq i \leq k$

$$\sup_{\mu \in \mathcal{K}_{k-i}} E(S^{(1)}_{\mu, 2}) = q^{-i} + q(1-q)^i / P^*$$

where $q = 1 - P^*$.

Proof.

$$\begin{aligned}
 & \sup_{\mathcal{E} \subset [k-j]} E(S^{t+1}_{\mathcal{E}}) \\
 &= \sup_{\mathcal{E} \subset [k-j]} \sum_{i=1}^j P(\text{select } r_i^{t+1} | \mathcal{E}) \\
 &\leq \sup_{\mathcal{E} \subset [k-j]} \sum_{i=1}^j P(\max_{r=1}^i r_i^{t+1} = d) \\
 &= \frac{j}{i+1} (1 - \frac{i}{r+1} E(-d)) \\
 &= j - \sum_{i=1}^j q^i \\
 &= j - q(1-q^i)/P^*
 \end{aligned}$$

where $q = (1-P^*)$.

Theorem 1, 3.24. $\sup_{\mathcal{E} \subset [k-j]} E(S^{t+1}_{\mathcal{E}})$ is increasing in j , hence

$$\sup_{\mathcal{E} \subset [k-j]} E(S^{t+1}_{\mathcal{E}}) \geq \sup_{\mathcal{E} = \emptyset} E(S^{t+1}_{\mathcal{E}}) = k - q(1-q^k)/P^*. \quad (1, 3.24)$$

Proof. Since the function

$$f(x) = x - ab^{x+1}$$

is increasing in x , for $0 < a < 1$, $0 < b < 1$, and $0 < x < \infty$.

In Case I of Gupta (1965) showed that

$$\sup_{\mathcal{E}} E(S^{t+1}_{\mathcal{E}}) = kp^*k. \quad (1, 3.24)$$

When the ordering prior about the unknowns is either unknown or known, we can use the selection procedure of Gustafsson and Broström (1975) where the ordering of the sample means as the ordering of unknowns is known, and apply the selection procedure which is currently used in the literature prior. With the latter approach, the substitution rule states that the isotonic regression of the sample means starts to the left of the true sample means, and that the selection procedures $\pi_{ij}^{(1)}, i = 1, 2, 3$, and $\pi_{ij}^{(2)}, j = 2, 3$ are the same as $\pi_{ij}^{(3)} (i = 1, 2, 3, 4)$, respectively, and the selection procedures $\pi_{ij}^{(1)}, i = 1, 2, 3, 4$ become $\pi_{ij}^{(4)}, i = 1, 2, 3$, respectively, which are equivalent to the procedures proposed by Gustafsson (1975) and Broström (1977), independent.

It has been proved in some quite general situation, and verified by using Monte Carlo technique in some selected cases by Gustafsson and Broström (1977), separately, that $\pi_p (\pi_{ij}^{(1)})$ is slightly better than π_4 . The values $d_{ijk}^{(j)}$ in the j th step of the procedure $\pi_{ij}^{(1)}$, $j = 1, 2, 3, 4$, are given by Broström (1977) as follows:

$$d_{ijk}^{(1)} = \frac{1}{\pi_{ij}^{(1)}(1 - (\beta^*)^j)}, \quad (1.1)$$

$$\int_{-\infty}^{\beta^*} \pi_{ij}^{(1)}(x + d_{ijk}^{(1)}) \phi(x) dx = \beta^*, \quad (1.2)$$

$$\int_0^{\beta^*} \pi_{ij}^{(1)}(xd_{ijk}^{(1)}) \phi(x) dx = \beta^*, \quad (1.3)$$

and

$$\int_{\beta^*}^{\beta^*} \int_{-\infty}^{\beta^*} \pi_{ij}^{(1)}(x + y)d_{ijk}^{(1)}(y) \phi(x) dy dx = \beta^*, \quad (1.4)$$

where ϕ is the density of $N(0, 1)$.

3.1. Selection Procedure for Scale Parameter of Gamma Population

Suppose we have $k+1$ independent populations π_1, \dots, π_{k+1} . The population π_i has a gamma density function

$$g(x; \alpha_i, \beta_i) = \frac{1}{\Gamma(\alpha_i)} \frac{x^{\alpha_i-1}}{\beta_i^{\alpha_i}} e^{-x/\beta_i}, \quad x > 0.$$

Example. The ordering prior of $\alpha_1, \dots, \alpha_k$ after β_1, \dots, β_k is given by $\alpha_1 < \dots < \alpha_k$. Note that the values of β_1, \dots, β_k and β_{k+1} are unknown, β_j 's are known.

In this section we define population π_i to be the i -th most scale population if the scale parameter $\beta_i < \beta_{k+1}$. Let $\alpha_i = \alpha_{i+k+1}$, then the parameter space is denoted by \mathbb{R}^k , where

$$\{x \in \mathbb{R}^{k+1} : \alpha_1 < \dots < \alpha_k, \beta_{k+1} = 0\}$$

is a subspace of $(k+1)$ -dimension Euclidean space \mathbb{R}^{k+1} .

Suppose we have independent observations x_{ij} ($i = 1, \dots, n_j$) from population π_i , ($i = 1, \dots, k$). Let $\bar{x}_{ij} = \bar{x}_{ij}/n_j$, then

$$x_{ij} - \frac{n_j}{j} \bar{x}_{ij}/n_j \text{ has density } g(x_{ij} - \bar{x}_{ij}/n_j),$$

and

$$\bar{x}_{ij}/n_j \text{ has density } g(\bar{x}_{ij}/n_j, 1/n_j).$$

Suppose our goal is to select a subset which contains all the populations under the ordering prior with probability greater than or equal to P_k , a predection test value between zero and one.

Let $\pi_{(1)}, \dots, \pi_{(k+1)}$ be the subsequence of population after π_{k+1} such that $\pi_{(j)} \in \Pi_{j-1}$ where

$$\begin{aligned} \hat{\gamma}_i &= \min(\hat{\gamma}_i, \hat{\gamma}_0 \wedge \hat{\gamma}_{i+1}), \quad \text{if } i = 1, \dots, m-1, \\ \hat{\gamma}_m &\in \{\hat{\gamma}_k \mid k = 0\}, \quad \text{if } i = m, \\ \hat{\gamma}_0 &\in \{\hat{\gamma}_0 \wedge \hat{\gamma}_1\}, \quad \text{if } i = 0. \end{aligned}$$

1.4.1. Proposed Selection Procedures $\hat{\gamma}_i$, $i = 6, 7, 8, 9$

Case I. Control γ_0 known and common sample size n .

Definition 1.4.1. The selection procedure $\hat{\gamma}_6$ is defined as follow.

Step 1. Select $\hat{\gamma}_i$, $i = k$ and stop, if

$$\hat{x}_{k:k} \leq c_{k:k} \wedge 0$$

otherwise reject $\hat{\gamma}_k$ and go to step 2.

Step 2. Select $\hat{\gamma}_i$, $i = k - 1$ and stop, if

$$\hat{x}_{k-1:k} \leq c_{k-1:k} \wedge 0$$

otherwise reject $\hat{\gamma}_{k-1}$ and go to step 3.

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Step k-1. Select $\hat{\gamma}_i$, $i = 2$ and stop, if

$$\hat{x}_{1:k} \leq c_{1:k} \wedge 0$$

otherwise reject $\hat{\gamma}_2$ and go to step k.

Step k. Select $\hat{\gamma}_1$ and stop, if

$$\hat{x}_{1:k} \leq c_{1:k} \wedge 0$$

otherwise reject $\hat{\gamma}_1$.

for each $i \in \{1, \dots, k\}$, i.e., to examine the smallest value $c_{1,i,k}$ such that the procedure δ_6 satisfies the B^* -condition.

Theorem 1.4.1. Assume we have random sample size n and $\rho_{ij} = \rho$ for all $i \in \{1, \dots, k\}$ and the constant $c_{1,i,k} (\leq 1)$ is determined by the condition

$$P(\theta_{ij} < c_{1,i,k}) \geq P^*, \quad i = 1, \dots, k, \quad (1.4.1)$$

where

$$U_i = \max_{j \leq s(i)} \left(\frac{Y_j + \epsilon_{ij}}{Y_j - \epsilon_{ij}} \right)^{\frac{1}{2}}, \quad (1.4.2)$$

and Y_i are i.i.d. with density

$$\pi(\cdot) \propto \frac{1}{n}, \text{ and cut-off } \beta(\cdot) \propto \frac{1}{n}$$

then the procedure δ_6 satisfies the B^* -condition.

Proof. For any i , $1 \leq i \leq k$, if the unknown true θ_{ij} ($j \neq i$) is fixed (i there are k good populations), then, under the procedure δ_6 ,

$$\begin{aligned} & \inf_{c \leq 1} P(\theta_{ij} < c) \\ &= \inf_{c \leq 1} P_{\lambda}(U_i < \frac{X_{ij} + \epsilon_{ij}}{X_{ij} - \epsilon_{ij}}) \\ &= \inf_{c \leq 1} P_{\lambda}\left(\frac{X_{ij} + \epsilon_{ij}}{X_{ij} - \epsilon_{ij}} < \frac{c + \epsilon_{ij}}{c - \epsilon_{ij}}\right) \\ &= \inf_{c \leq 1} P_{\lambda}\left(\frac{(c + \epsilon_{ij})(X_{ij} - \epsilon_{ij})}{(c - \epsilon_{ij})(X_{ij} + \epsilon_{ij})} < 1\right) \\ &= \inf_{c \leq 1} P_{\lambda}\left(\frac{(c + \epsilon_{ij})(X_{ij} - \epsilon_{ij})}{(c - \epsilon_{ij})(X_{ij} + \epsilon_{ij})} < 1\right) \\ &= \inf_{c \leq 1} P_{\lambda}\left(\frac{(c + \epsilon_{ij})(X_{ij} - \epsilon_{ij})}{(c - \epsilon_{ij})(X_{ij} + \epsilon_{ij})} < 1\right) \\ &= \inf_{c \leq 1} P_{\lambda}\left(\frac{(c + \epsilon_{ij})(X_{ij} - \epsilon_{ij})}{(c - \epsilon_{ij})(X_{ij} + \epsilon_{ij})} < 1\right) \end{aligned}$$

$$\begin{aligned}
&= \inf_{\mathbf{c}_{i:k}} P_{\mu}(\cup_{j=i}^k (\max_{1 \leq s \leq j} \min_{s \leq t \leq k} \frac{Y_{s:s} + \dots + Y_{t:t}}{t-s+1} \leq c_{i:k})) \\
&= P_{\mu \otimes \nu^k}(\cup_{j=i}^k (\max_{1 \leq s \leq j} \min_{s \leq t \leq i} \frac{Y_{s:s} + \dots + Y_{t:t}}{t-s+1} \leq c_{i:k})) \\
&= P(\max_{1 \leq s \leq i} \frac{Y_{s:s} + \dots + Y_{i:i}}{i-s+1} \leq c_{i:k}) \\
&= P(U_i \leq c_{i:k}),
\end{aligned}$$

where Y_i 's are i.i.d with density $a(\cdot; \nu, \frac{1}{n})$,

$$\mathbf{v}^* = (\underbrace{0, 0, \dots, 0}_{i+1}, \dots, \mathbf{v})$$

and

$$U_i = \max_{1 \leq s \leq i} \frac{Y_{s:s} + \dots + Y_{i:i}}{i-s+1}.$$

Corollary 1.4.1. $c_{i:k} = c_{i:i}$, $i = 1, \dots, k$.

For any $x > 0$, let $S_n = \sum_{i=1}^n (Y_i - x)$, $n = 1, 2, \dots$, $\nu_0 = 0$. Then
 $0 < P(Y_i - x > 0) < a(x, \nu, \frac{1}{n}) < 1$, the distribution of $Y_i - x$ is concentrated on a half-axis. By Theorem 1.4.1, the probability generating function of cumulative distribution functions $P(U_k > x)$, $k = 1, 2, \dots$, is given by

$$\exp \left[\frac{x}{1-\nu} + b(kn, \nu, 1) \right].$$

Hence by Theorem 1.3.6, we have the following recurrence formula for all $x > 0$:

$$P(U_{k+1} = x)$$

$$= \sum_{j=0}^{k+1} P(U_{k+1} = x) + G((j+1)n; (j+1), 1) \quad (1.4.4)$$

where

$$P(U_0 = x) = 1.$$

When $x = 0$, both sides of Equation (1.4.4) equal to zero, hence it also holds for $x = 0$.

Note that

$$\begin{aligned} P\left(\frac{1}{r} \sum_{i=1}^r Y_i = x\right) &= G(x; r, \frac{1}{rn}) \\ &= G(xn; r, 1). \end{aligned} \quad (1.4.5)$$

The values $c_{k:k}(P^*, r, n)$ which satisfy Equation (1.4.1) are tabulated in Table IV for $k = 1(1)10$, $P^* = .29, .45, .60, .75$, $r = 2, 4, 6$, and $n = 4, 6, 8, 10, 15, 20$.

Lemma 1.4.1. $c_{ik} = c_{i+1:k}$ for all $1 \leq i \leq k-1$.

Proof. The constants c_{ik} ($i = 1, \dots, k$) are determined by (1.4.1), respectively,

$$U_{i+1} - U_i \geq 0 \text{ implies } c_{ik} = c_{i+1:k} \text{ for all } 1 \leq i \leq k-1.$$

Theorem 1.4.7. The selection procedure φ will not be changed if the monotone estimators \hat{Y}_{ijk} , $i = 1, \dots, k$ are replaced by $\hat{c}_{ijk} \hat{Y}_{ijk} + \hat{c}_{i+1:k} \hat{Y}_{i+1:k}$.

where

$$\begin{aligned} Y_{j+1} &= \max_{1 \leq i \leq j} T_{ij} \\ &= \max_{\substack{1 \leq s \leq j \\ 1 \leq t \leq k}} \frac{T_{st}}{\min_{1 \leq u \leq k} T_{su}} \end{aligned}$$

Proof. The proof is similar to that of Theorem 1.4.1.

Next, we define a selection procedure by using an efficient estimator and a fixed constant which depends on P^* , k , sample size n and constant c_3 .

Definition 1.4.2. The selection procedure γ_j is defined by

γ_j : Select x_j if and only if $X_{j+1} < c_0$ ($j = 1, \dots, n$)

where $c_0 \in (1)$ is the smallest value such that procedure γ_n satisfies the P^* -condition.

Corollary 1.4.2. The constant $c_0 \in (P^*, k, n)$ of the selection procedure γ_j equals to $c_{k;k}$ which is determined by the following rule.

Proof. Follows immediately from Theorem 1.4.1 and (1.4.1).

Definition 1.4.3. The selection procedure γ is defined by the following steps:

Step 1: Select x_1 if and only if $X_2 < c_0$
 \vdots
 \vdots
 \vdots
 \vdots

otherwise proceed to Step 2

Step 2. Select c_j , $j = k+1$, and stop, if

$$\begin{cases} x_{k+1} < c_{k+1} \\ c_{k+1} > 0 \end{cases}$$

otherwise reject c_{k+1} and go to step 1.

Step 3.1. Select c_j , $j = 1$, and stop, if

$$\begin{cases} x_1 < c_1 \\ c_1 > 0 \end{cases}$$

otherwise reject c_1 and go to step k.

Step k. Select c_1 and stop, if

$$\begin{cases} c_1 < 0 \\ c_1 > 1 \end{cases}$$

otherwise reject c_1 .

where the c_j 's are the smallest real values (> 1) such that the $(x_j - c_j)_+$ satisfies the P*-condition.

where $\{x_i\}_{i=1}^n$ the i -th initial point are determined by

$$x_i = \frac{1 + (-1)^{i+1}}{2} c_i, \quad i = 1, \dots, n, \quad \text{and}$$

and the efficiency is defined by $\eta = \frac{c_1}{c_{k+1}}$.

REFERENCES

[1]

$$= \inf_{\tau_{i+1}} P_i(\tau_i) \left(\frac{\tau_i}{\tau_{i+1}} - c_{i+1} \right)^{p_i}$$

$$= \inf_{\tau_{i+1}} P_i(\tau_i) \left(\frac{\tau_i}{\tau_{i+1}} - c_{i+1} \right)^{p_i}$$

$$= P_i(\tau_i) \left(\frac{\tau_i}{\tau_{i+1}} - c_{i+1} \right)^{p_i}$$

$$= P(Z_i < c_{i+1})$$

where $Z_i = \frac{\tau_i}{\tau_{i+1}}$, $i = 1, \dots, k$ are i.i.d. with the same density.

$g(\cdot; \tau_i, 1), \tau^* = (\underbrace{\tau_0, \dots, \tau_0}_{i+1}, \dots, \tau_k)$. Hence c_j , $j = 1, \dots, m$ is determined by (1.4.6). If $\tau_1 < \dots < \tau_k$, then $c_1 < \dots < c_k$.

The following selection procedure π_0 was given by Gupta and Edel (1958).

Definition 1.4.4. The selection procedure π_0 is defined by

(a) Select τ_i if and only if $\frac{\tau_i}{\tau_{i+1}} > c_i$, $i = 1, \dots, k$
where c_i is determined by

$$\frac{k-i+1}{i+1} \int_{c_i}^{c_i} \frac{e^{-\lambda u}}{u} du = \alpha^* \quad (1.4.7)$$

for $c_i > 0$, $i = 1, \dots, k$, it turns to

$$\frac{1}{\lambda} \int_{c_i}^{c_i} \frac{e^{-\lambda u}}{u} du = \alpha^* \quad (1.4.8)$$

the left-hand side is the c.d.f. of $\frac{1}{2} \chi^2$ with r degrees of freedom, hence the value c_1 can be easily solved with the help of a table of chi-square distribution.

Application to the Selection of Variance of Normal Population

If $\sigma_{ij}^2 = \sigma_i^2$, $i = 0, 1, \dots, k$ are the scale parameters for the $k+1$ normal populations and y_{ij} ($j = 1, \dots, n$; $i = 1, \dots, k$) are the observations on the population γ_i with the mean μ_i (known), we assume that the order $\sigma_1^2 < \dots < \sigma_k^2$ is known.

In the application of selection procedure ϵ or γ_2 , what we need to do is to evaluate the isotonic regression of S_i^2 which is the sample variance of population γ_i , $i = 1, \dots, k$ and denote it by S_{i+k}^2 , $i = 1, \dots, k$, then directly apply ϵ_6 or γ_2 . The constant we need is determined by Equations (1.4.2) and (1.4.4) where we replace n by n , the reason being that (nS_i^2/σ_i^2) has χ^2_n distribution with n degrees of freedom and $t_{ij} = S_{i+k}^2/\sigma_i^2$ has the c.d.f.

$$G(2nt; n, \beta) = G(t; n, \frac{1}{n}),$$

hence

$$P\left(\frac{1}{n} \sum_{i=1}^r y_{ij} - t\right) = G(trn; rn, 1).$$

The application of γ_1 is similar to that of ϵ_6 (see Indiana University (1969)). What we need to do is to replace y_{ij} in γ_1 and ϵ_6 by $y_{ij} - t$, replace n in Equation (1.4.6) and (1.4.7) by n , $i = 1, \dots, k$.

Remark 1.4.1. γ_0 (Gupta and Sobel (1958)) does not depend on the underlying prior and the sample sizes for each population need not be equal.

If the means μ_i , $i = 1, \dots, k$ are unknown and common, and if we let $n = 1$, let $S_i^{(2)} = \sum_{j=1}^n (\bar{X}_{ij} - \bar{X}_i)^2/n-1$ and use $n = 1$ in place of n in (1.4.4), (1.4.6) and (1.4.8) which determined the constants $c_{i;k}$, c and c' for γ_6 , γ_8 and γ_9 , respectively.

1.4.2. Selection Procedure $\gamma_i^{(2)}$, $i = 6, 7, 8, 9$

Case II. γ_0 unknown.

The assumptions are the same as in Case I except that n_0 observations, viz., x_{01}, \dots, x_{0n_0} are taken on γ_0 .

For selection procedure $\gamma_6^{(2)}$, the inequalities defining the procedure and corresponding to $\bar{X}_{i;k} - c_{i;k} \gamma_0$ ($i = 1, \dots, k$) are replaced by $\bar{X}_{i;k} - c_{i;k}^{(2)} \gamma_0$ ($i = 1, \dots, k$), respectively. The equation determining $c_{i;k}^{(2)}$ is obtained as before and is given by

$$\int_0^\infty P(U_i - c_{i;k}^{(2)} > t) dt = 1,$$

where $P(U_i > t)$ is the survival function in Theorem 1.4.1, i.e., the r.d.f. of γ_0 if population γ_0 .

If population γ_0 has gamma distribution with density $\alpha \gamma_0^{\alpha-1} e^{-\gamma_0}$ (α_0 known and γ_0 unknown), then $P(U_i > t) = n_0 \Gamma(t/n_0)$.

For selection procedure $\gamma_7^{(2)}$, the inequality defining the procedure is,

$$x_{k+1} < c^* x_0$$

and it can be shown $c^* = \frac{c}{k+1}$.

For selection procedure $\frac{(k)}{g}$, the inequality defining the prior estimate and corresponding the $(k-i)$ th step is

$$\frac{x_i - x_0}{x_i} < \frac{\lambda_0}{\lambda_0}, \text{ where } \lambda_0 = n_0/\alpha.$$

The equation determining c of $\frac{(k)}{g}$ is given by

$$\int_0^{t_{k-i}} \left[\frac{c - u^2}{n} \right] \frac{e^{-u}}{u} du \int_{t_{k-i}}^{t_k} \frac{e^{-t}}{t^2} dt = p^*, \quad (1.1.1)$$

For selection procedure $\frac{(k)}{g}$, the inequality defining the prior estimate is

$$\frac{x_i - x_0}{x_i} < \frac{\lambda_0}{\lambda_0}, \quad (1.1.2)$$

and the equation determining c is given by Gupta and Soper (1970) as follows:

$$\int_0^{t_{k-i}} \left[\frac{c - u^2}{n} \right] \frac{e^{-u}}{u} du \int_{t_{k-i}}^{t_k} \frac{e^{-t}}{t^2} dt = p^*, \quad (1.1.3)$$

1.5. Selection Rules for the Location Parameter (continued)

Ordering Prior Assumption

Suppose that we have only a partial ordering information about the location parameters, that is, type I or type II ordering.

\mathbb{R}^k is a space for a set of k elements.

For a given k , let \mathcal{P}_k be the set of all partitions of \mathbb{R}^k .

Let $\mathcal{B}_k = \{B_1, \dots, B_m\}$ be a partition of \mathbb{R}^k into m disjoint nonempty sets, say B_1, \dots, B_m so that $x_i \in B_j \iff x_i \in \mathbb{R}^j$,

and for each $b \in B_1, \dots, B_m$ there is a unique $i(b)$ such that $b \in \mathbb{R}^{i(b)}$.
There is no order relation among the elements of B_1, \dots, B_m .

Let $b_{ij} \in B_{i(j)}$, the number of elements in $B_{i(j)}$,
so we have

$$\frac{b_{ij}}{b_{i(j)}} < 1$$

If we denote the new parameter by $t_{ij} = b_{ij}/b_{i(j)}$, then t_{ij} is a parameter which is < 1 . We can now complete \mathcal{P}_k to \mathcal{P}_k' with an induced partial order.

Example. Suppose $k = 3$, and we have a partial ordering on \mathbb{R}^3 given by
 $1 \leq 2 \leq 3$, and $a \leq c$ if $a \leq b \leq c$. Then t_{ij} represent this partial ordering, as follows:



Figure 1. An initial partial ordering.

Then we have an induced partial order from $\{x_1, x_2, \dots, x_n\}$ to $\{y_1, y_2, \dots, y_n\}$ as in Figure 2.

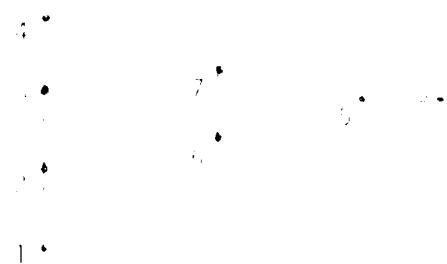


Figure 2. Induced partial ordering.

Now

$$B_0 = \{x_1^*, x_2^*\}$$

$$B_1 = \{x_1^*, x_2^*, x_3^*\}$$

$$B_2 = \{x_1^*, x_2^*, x_3^*, x_4^*\}$$

It is clear that the induced partial order is not unique. For example, we can partition $\{x_1, x_2, \dots, x_n\}$ into three other subsets B_0^1, B_1^1, B_2^1 , where

$$B_0^1 = \{x_1^*, x_2^*\}$$

$$B_1^1 = \{x_1^*, x_2^*, x_3^*, x_4^*\}$$

$$B_2^1 = \{x_5^*, x_6^*, x_7^*, x_8^*\}$$

From the local to global ordering, a selection procedure $\tilde{\sigma}_n$ is defined as follows:

Definition 1. Let us define a selection procedure $\tilde{\sigma}_n$ as follows:

Suppose $x_{1,2}, \dots, x_n$ are the incident points and that the elements of $x_{1,2}, \dots, x_n$ satisfy the following properties: (i) the incident property

selection procedure — for each subset \mathcal{S}_j we have to find the maximum value

of the function $\text{int}^P(\mathcal{S}_j)$ and the probability of a correct selection is not less than $1 - \alpha_j^P$.

For the selection procedure of the best subset \mathcal{S}_j from the set \mathcal{S} of all subsets \mathcal{S}_j we may use selection procedure $\mathcal{S}_{\text{opt}}^P$ or $\mathcal{S}_{\text{opt}}^{P*}$ with $\alpha_j^P = \alpha_j^{P*}$.

Remark 1.5.1. The selection procedure $\mathcal{S}_{\text{opt}}^P$ or $\mathcal{S}_{\text{opt}}^{P*}$ can be used in the following cases:

(a) if

$$\text{int}^P(\mathcal{S}) > 0$$

$$\text{int}^P(\mathcal{S}) \geq \text{int}^P(\mathcal{S}_j)$$

$$\text{int}^P(\mathcal{S}) > \text{int}^P(\mathcal{S}_j)$$

$$\mathcal{S} \in \mathcal{S}_j$$

$$\begin{aligned} &\text{int}^P(\mathcal{S}) > \text{int}^P(\mathcal{S}_j) \\ &\mathcal{S} \in \mathcal{S}_j \\ &\mathcal{S} \neq \mathcal{S}_j \end{aligned}$$

where α_j^P is the parameter of accuracy associated with the subset \mathcal{S}_j .

Remark 1.5.1. For the selection problem of the best subset \mathcal{S}_j from the general distributions, when α_j^P is not fixed, the following algorithm for the selection procedure can be applied: scanning the subsets \mathcal{S}_j in the order

1.6. Computation of the Bayesian procedure coefficients

1.6.1. The location parameters of normal distribution

In this section we use Monte Carlo technique to estimate the performance of the selection procedures $\mathcal{S}_{\text{opt}}^P$ and $\mathcal{S}_{\text{opt}}^{P*}$ with $\alpha_j^P = \alpha_j^{P*}$.

K independent populations, each population with distribution $N(\mu_i, \sigma^2)$, with common known variance σ^2 and common sample size n_i . Assume that the mean μ_0 of the control is known; without loss of generality we assume that $\mu_0 = 0$ and $\sigma/\sqrt{n} = 1$.

In the simulation, we use Rubin and Trinkle's PVP-Param Variable Package, Purdue University Computing Center, to generate random numbers. For each k , we generated one random number (variable) for each population, then applied each selection procedure separately and repeated it ten thousand times; we used the relative frequency as an approximation of the exact values of the associated performance characteristics for each procedure. In Table V we use the following notations:

$\mu = (\mu_1, \dots, \mu_k)$, μ_i is the parameter of population i ,

PS = $P(CS)$

PI = $P(\text{correctly rejecting all bad populations})$

PC = $P(\text{correct classification of all populations})$

where the correct classification means that we select all good populations and reject all bad populations.

EE = Expected number (size) of bad populations contained in the selected subset,

$$\sum_{\substack{i=1 \\ i \neq 0}}^K \left(\mu_i - \mu_0 \right)^2 p_i^2 t_i \quad (\mu_i \text{ is selected})$$

ES = Expected size of the selected subset,

$k = 3, 4, 5, 6$. For each value of k , we assume that the true population is the one and only one parameterization with performance that is less than the control ($\alpha = 0.05$). A glance at the table in Section 3 shows that the performance can roughly be ordered as follows:

• • • • •

The procedure τ_3 is the best one performed by both methods. τ_2 and τ_1 are very close and both are better than τ_4 . The choice is based on the characteristics of the problem. As mentioned in Section 4.1, if λ is small, we can choose τ_1 or τ_2 . If λ is large, we can choose τ_3 or τ_4 . As the number of model components increases, the value of the parameter λ , α_1 and β_1 respectively, we find that τ_1 , τ_2 , τ_3 and τ_4 ($i = 1, 2, 3, 4$) is reduced, α_1 and β_1 mostly increase. This means that when k is increasing and the additional population, τ_1 and τ_2 are the best one for the model estimation.

Table V contains the detailed information about the performance characteristics of the two types of traffic in the network. It can be observed from the table that we have two heavy traffic flows in our network. The first one is the performance characteristic of the heavy traffic flow in the company office premises.

In Table 7, we compare the results obtained by using the two different methods. It is evident that both the methods give similar results. The values of the error function in Table 7 are the same as those given in Table 4.

population, then the \hat{P}_k and \hat{P}_{k+1} are identical. This is because \hat{P}_k is the best estimate of the population given the data up to time k .

In Table V.4 we compare the memory required to store the population by both methods. The amount of memory required and that performance is excellent. We also show in Table V.5 the same results as before.

In Table V.6 we compare the performance of the two procedures. For the first procedure, the best estimate of the population P_k , P_{k+1} and \hat{P}_{k+1} are given for each value of ρ . Again, \hat{P}_{k+1} is the best. If we compare \hat{P}_{k+1} and P_{k+1} it is clear that \hat{P}_{k+1} does not change the small difference between P_{k+1} and P_k . The fractions $\lambda_1 P_{k+1}^2$ and $\lambda_2 P_{k+1}$ increase with ρ . This implies from equation (5) that the performance of the second procedure \hat{P}_{k+1} is determined by the fraction $\lambda_2 P_{k+1}$. In fact, $\lambda_2 P_{k+1}$ increases as ρ increases. From these five tables, it is apparent that the best estimate of the population P_k is given by \hat{P}_{k+1} . It is also apparent that the second procedure is better than the first if ρ is correctly estimated or at least close to the true value. It is also better than the other two procedures if ρ is not estimated correctly, otherwise the other two procedures are better.

Finally, we compare the performance of the two procedures for different values of ρ .

In this section, we assume the \hat{P}_{k+1} to be the best estimate of the population P_{k+1} .

population has exponential distribution $\text{Exp}(\lambda)$ with scale parameter λ and with unknown scale parameter λ (see, e.g., [1]).

In the simulation study, the estimator $\hat{\gamma}_g$ is compared with $\hat{\gamma}_P$ and $\hat{\gamma}_R$, where we used Rubin and Stakelum's PIP and estimator $\hat{\gamma}_R$ was proposed by Rubin and Stakelum [1]. The population λ is simulated from $\text{Exp}(1)$ with specified exponential distribution function. The simulation results are based on two thousand sets of random samples. The simulation results are given in Table VI.1 and Table VI.2, respectively, for $\lambda = 1$ and $\lambda = 1.6$.¹ but the results of a bad population function are also given for $\lambda = 1$ and $\lambda = 1.6$.¹ A bad population has a parameter which is generated from uniform distribution (assume one).

A good population means that the parameter is 1.00. The ratio of $\hat{\gamma}_g$ to $\hat{\gamma}_P$ is close to one. The results of Table VI.1 and Table VI.2, and especially Table VI.2, show that we have the performance $\hat{\gamma}_g \approx \hat{\gamma}_P < \hat{\gamma}_R$ in the aspect of efficiency, i.e., that $\hat{\gamma}_g$ is slightly better than $\hat{\gamma}_P$. There is only one bad situation in

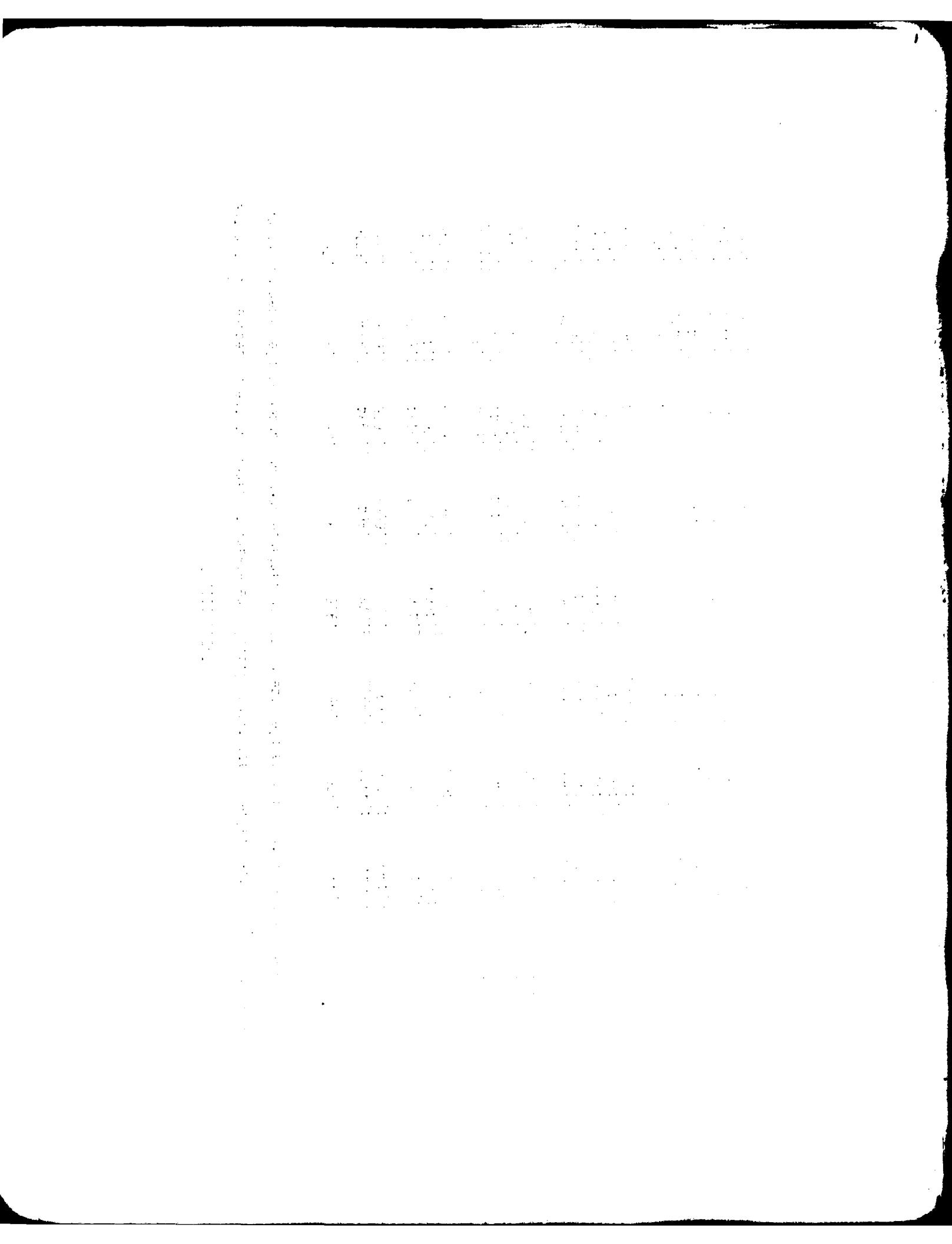
Table of $d_{1,k}^{(1)}$ values, calculated by the direct and the indirect methods, to carry out the procedure π_1 for the normalized quadratic loss function ordering prior.

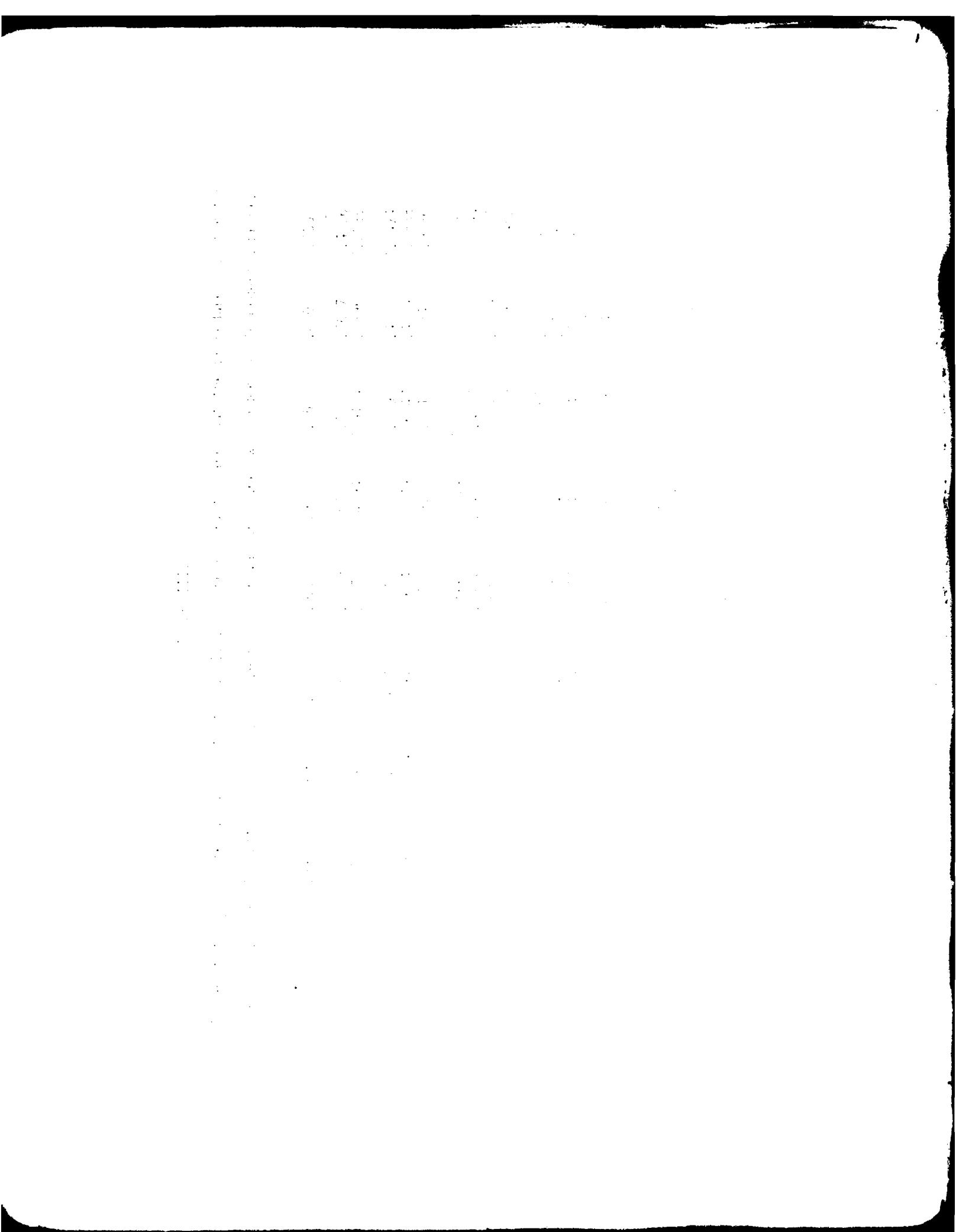
k	$d_{1,k}^{(1)}$	$d_{1,k}^{(2)}$	$d_{1,k}^{(3)}$	$d_{1,k}^{(4)}$	$d_{1,k}^{(5)}$	
1	2.3764	1.6666	1.4480	1.4346	1.4346	
2	2.3347	1.6756	1.6200	1.3876	1.4444	
3	2.3329	1.6767	1.6447	1.4986	1.5555	
4	--	--	1.6515	1.4686	1.5555	
5	--	--	1.6536	1.4966	1.5555	
6	--	--	--	--	1.5555	
7	1.3400	1.9787	1.6574	1.4946	1.5555	
$d_{1,k}^{(1)}$	$d_{1,k}^{(2)}$	$d_{1,k}^{(3)}$	$d_{1,k}^{(4)}$	$d_{1,k}^{(5)}$		
1	1.36	1.60	1.31	1.10	1.10	
2	1.0664	1.6466	1.6356	1.6366	1.6366	
3	1.1736	1.6666	1.6666	1.6666	1.6666	
4	1.1436	1.6767	1.6767	1.6767	1.6767	
5	1.1426	1.6767	1.6767	1.6767	1.6767	
6	1.1426	1.6767	1.6767	1.6767	1.6767	
7	1.1426	1.6767	1.6767	1.6767	1.6767	

For the first four values of k , the values of $d_{1,k}^{(1)}$ were calculated by the direct method, while the remaining values were calculated by the indirect method.

Table of $\alpha_{\text{per}}^{(2)}$ Additive correction to the
procedure $\frac{\ell_1}{\ell_2}$ for the normal measurement problem

$\alpha_{\text{per}}^{(2)}$	ℓ_1	ℓ_2	$\ell_1 + \ell_2$	$\ell_1 - \ell_2$
0.00	0.0000	0.0000	0.0000	0.0000
0.05	-0.0000	-0.0000	-0.0000	-0.0000
0.10	-0.0000	-0.0000	-0.0000	-0.0000
0.15	-0.0000	-0.0000	-0.0000	-0.0000
0.20	-0.0000	-0.0000	-0.0000	-0.0000
0.25	-0.0000	-0.0000	-0.0000	-0.0000
0.30	-0.0000	-0.0000	-0.0000	-0.0000
0.35	-0.0000	-0.0000	-0.0000	-0.0000
0.40	-0.0000	-0.0000	-0.0000	-0.0000
0.45	-0.0000	-0.0000	-0.0000	-0.0000
0.50	-0.0000	-0.0000	-0.0000	-0.0000
0.55	-0.0000	-0.0000	-0.0000	-0.0000
0.60	-0.0000	-0.0000	-0.0000	-0.0000
0.65	-0.0000	-0.0000	-0.0000	-0.0000
0.70	-0.0000	-0.0000	-0.0000	-0.0000
0.75	-0.0000	-0.0000	-0.0000	-0.0000
0.80	-0.0000	-0.0000	-0.0000	-0.0000
0.85	-0.0000	-0.0000	-0.0000	-0.0000
0.90	-0.0000	-0.0000	-0.0000	-0.0000
0.95	-0.0000	-0.0000	-0.0000	-0.0000
1.00	-0.0000	-0.0000	-0.0000	-0.0000





1. The first step in the process of determining the nature of the offense is to identify the offense itself. This is done by examining the facts of the case and determining what specific offense has been committed. This may involve reviewing witness statements, physical evidence, and other relevant information.

2. Once the offense has been identified, the next step is to determine the appropriate legal classification for the offense. This involves considering factors such as the severity of the offense, the intent of the offender, and the circumstances surrounding the offense. This may involve consulting legal codes and statutes, as well as seeking advice from legal professionals.

3. After the offense has been identified and classified, the next step is to determine the appropriate punishment or sentence for the offender. This involves considering factors such as the gravity of the offense, the offender's criminal history, and the impact of the offense on the victim. This may involve consulting legal codes and statutes, as well as seeking advice from legal professionals.

4. Finally, the last step in the process of determining the nature of the offense is to implement the sentence or punishment. This may involve sentencing the offender to jail, probation, or fines, or taking other legal action. It may also involve providing rehabilitation or treatment services to the offender.

the value $\rho_{\text{max}} = 0.8$ in the first column of Table 1, we can see that the error rate π_0 for the quantile-based procedure is within 0.05 of the nominal significance level under the null hypothesis.

TABLE 1

	$\alpha = 0.05$	$\alpha = 0.01$	$\alpha = 0.001$	$\alpha = 0.0001$
1	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
2	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
3	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
4	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
5	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
6	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
7	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
8	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
9	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
10	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
11	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
12	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
13	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
14	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
15	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
16	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
17	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
18	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
19	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000
20	0.0513 0.0512	0.0100 0.0100	0.0009 0.0009	0.0000 0.0000

TABLE 1. Empirical Type I error rates for the quantile-based procedure under the null hypothesis.

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TABLE II (continued)

n₁(1)

	n ₁	n ₂	n ₃	n ₄	n ₅	n ₆	n ₇
1	1.6702	1.5958	1.5713	1.4976	1.4441	1.3717	
2	1.7039	1.6668	1.6367	1.4596	1.3980	1.3277	
3	1.7406	1.6634	1.6147	1.4561	1.3728	1.3037	
4	1.7440	1.6976	1.6157	1.4579	1.3711	1.3034	
5	1.7451	1.6895	1.6156	1.4584	1.3704	1.3031	
6	1.7455	1.6888	1.6157	1.4585	1.3707	1.3031	
7	1.7456	1.6879	1.6157	1.4586	--	--	
8	1.7457	--	--	--	--	--	
10	--	--	--	--	--	--	
1	2.79437	2.7567	2.5615	2.4395	2.4799	2.4449	
2	3.0126	2.5164	2.3951	2.6179	2.5136	2.4433	
3	3.0264	2.5223	2.7143	2.6396	2.5144	2.4434	
4	3.0301	2.8320	2.7156	2.6368	2.5156	2.4443	
5	3.0313	2.8328	2.7162	2.6374	2.5158	2.4443	
6	3.0317	2.8331	2.7164	2.6376	2.5164	2.4444	
7	3.0318	2.8331	2.7165	2.6377	--	2.4444	
8	3.0319	--	--	--	--	--	
10	--	--	--	--	--	--	
1	4.1496	3.1134	3.1304	3.1210	3.1258	3.1317	
2	4.1736	3.1268	3.1363	3.1636	3.1496	3.1398	
3	4.1744	4.1064	3.1865	3.1717	3.1569	3.1366	
4	4.17470	4.1092	3.1865	3.1737	3.1574	3.1413	
5	4.17493	4.1101	3.1866	3.1734	3.1577	3.1413	
6	4.17496	4.1114	3.1867	3.1736	3.1577	3.1413	
7	4.17498	4.1114	3.1867	3.1737	--	3.1413	
10	--	--	--	--	--	--	

the first time, and the first time I have seen it. It is a very large tree, and has a very large trunk.

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TABLE II
Calculated values of the energy and the constant terms of various states of the protonium-like atom formed by the interaction of the proton and the electron. The units of energy and constant term are MeV.

	E_p	E_e	C_p	C_e
$k = 4, \quad (-1, 0, 0)$				
P ₀	-1.9454	1.9454	-1.9454	1.9454
P ₁	-1.3894	1.3894	-1.3894	1.3894
P ₂	-1.3307	1.3307	-1.3307	1.3307
P ₃	-1.6146	1.6146	-1.6146	1.6146
P ₄	-1.6146	1.6146	-1.6146	1.6146
P ₅	-1.5599	1.5599	-1.5599	1.5599
$k = 4, \quad (-1, 0, 1)$				
P ₀	-1.9611	1.9611	-1.9611	1.9611
P ₁	-1.3744	1.3744	-1.3744	1.3744
P ₂	-1.3272	1.3272	-1.3272	1.3272
P ₃	-1.6259	1.6259	-1.6259	1.6259
P ₄	-1.6259	1.6259	-1.6259	1.6259
P ₅	-1.5721	1.5721	-1.5721	1.5721
$k = 4, \quad (-1, 0, 1, 0)$				
P ₀	-1.9630	1.9630	-1.9630	1.9630
P ₁	-1.3664	1.3664	-1.3664	1.3664
P ₂	-1.3144	1.3144	-1.3144	1.3144
P ₃	-1.6136	1.6136	-1.6136	1.6136
P ₄	-1.6136	1.6136	-1.6136	1.6136
P ₅	-1.5600	1.5600	-1.5600	1.5600
$k = 4, \quad (-1, 0, 1, 1)$				
P ₀	-1.9634	1.9634	-1.9634	1.9634
P ₁	-1.3715	1.3715	-1.3715	1.3715
P ₂	-1.3241	1.3241	-1.3241	1.3241
P ₃	-1.6127	1.6127	-1.6127	1.6127
P ₄	-1.6127	1.6127	-1.6127	1.6127
P ₅	-1.5611	1.5611	-1.5611	1.5611

simulation results for the optimal policies under different prior distributions and different values of the parameter L (Eq. (1) per unit of production).

The simulation results are presented in Table I.

Table I shows that the optimal policy is characterized by

(a) a low value of L ($L = 10$) and a large value of α ($\alpha = 0.95$);
(b) a small value of L ($L = 10$) and a small value of α ($\alpha = 0.90$);
(c) a large value of L ($L = 100$) and a small value of α ($\alpha = 0.90$);
(d) a large value of L ($L = 100$) and a large value of α ($\alpha = 0.95$).

Thus, the optimal policy is characterized by a small value of L and a large value of α .

Table II presents the simulation results for the optimal policy under different prior distributions and different values of the parameter L (Eq. (1) per unit of production).

The simulation results are presented in Table II.

Table II shows that the optimal policy is characterized by

(a) a small value of L ($L = 10$) and a large value of α ($\alpha = 0.95$);
(b) a small value of L ($L = 10$) and a small value of α ($\alpha = 0.90$);
(c) a large value of L ($L = 100$) and a small value of α ($\alpha = 0.90$);
(d) a large value of L ($L = 100$) and a large value of α ($\alpha = 0.95$).

Thus, the optimal policy is characterized by a small value of L and a large value of α .

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Thus, the optimal policy is characterized by a small value of L and a large value of α .

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Simulation results for the comparative performance of various selection procedures for the normal means under distributional conditions 1, 6, 11 under simple random sampling.

Table V, 3

Simulation results for the comparative performance of various selection procedures for the normal means problem (notations explained in Section 1.6.1) under simple ordering prior.

$\mu^* = .900$

$k = 2, \pi = (-1, +1)$

	1	2	3	4
P5	1.0000	1.0000	1.0000	1.0000
PI	.5405	.5343	.5337	.5272
PC	.5405	.5343	.5337	.5177
E1	.8331	.8337	.8351	.8394
E3	.77116	.77349	.74332	.73773
E5	.8331	.8337	.8351	.8394

$k = 3, \pi = (-1, +1, +1)$

	1	2	3	4
P5	.9957	.9957	.9957	.9970
PI	.5365	.5343	.5337	.5197
PC	.5397	.5343	.5337	.5197
E1	.8347	.8363	.8316	.8394
E3	.7759	.77349	.74155	.73919
E5	.8279	.8330	.8373	.8394

$k = 4, \pi = (-1, +1, +1, +1)$

	1	2	3	4
P5	.9924	.9957	.9957	.9970
PI	.5274	.5343	.5334	.5197
PC	.5197	.5343	.5337	.5197
E1	.8403	.8363	.8316	.8377
E3	.7765	.77349	.74155	.73919
E5	.8390	.8330	.8373	.8394

$k = 5, \pi = (-1, +1, +1, +1, +1)$

	1	2	3	4
P5	.9901	.9957	.9957	.9970
PI	.5217	.5343	.5337	.5197
PC	.5177	.5343	.5337	.5197
E1	.8461	.8341	.8316	.8377
E3	.7761	.77349	.74155	.73919
E5	.8341	.8330	.8373	.8394

TABLE VI. 1

Simulation results for the comparative performance of various statistical procedures for the gamma means problem (notation explained in section 1.6.1) under simple ordering constraint.

P^* = .500

$\kappa = 4, \mu = (.3,.6,.9,1.2)$

	α	β	γ	δ
PS	.0840	.0848	.0850	.0850
PI	.7629	.7745	.7752	.7752
PC	.1860	.1867	.1870	.1870
FI	.7980	.7995	.8005	.8005
ED	.0319	.0320	.0320	.0320
EF	2.7320	3.8190	3.8410	3.8410

$\kappa = 6, \mu = (.3,.6,.9,1.7,1.8)$

	α	β	γ	δ
PS	.0905	.0910	.0915	.0915
PI	.6990	.7090	.7045	.7045
PC	.0695	.0750	.0645	.0645
FI	1.5040	1.4560	1.5870	1.5870
ED	.1868	.1969	.0000	.0000
EF	4.4945	4.5470	4.5865	4.5865

$\kappa = 4, \mu = (.3,1,1.5,1.7)$

	α	β	γ	δ
PS	.0625	.0628	.0625	.0625
PI	.7625	.7667	.7670	.7670
PC	.2370	.2373	.2375	.2375
FI	1.0790	1.1110	1.1100	1.1100
ED	.5304	.5310	.0000	.0000
EF	2.2475	2.2710	2.2710	2.2710

$\kappa = 6, \mu = 1,1.5,1.7,1.8$

	α	β	γ	δ
PS	.0600	.0600	.0600	.0600
PI	.7600	.7600	.7600	.7600
PC	.1110	.1110	.1110	.1110
FI	1.0710	1.0710	1.0710	1.0710
ED	.5300	.5300	.0000	.0000
EF	2.2400	2.2400	2.2400	2.2400

TABLE VI.

Simulation results for the comparative performance of various selection procedures for the gamma mean problem (notation explained in section 1.6.1) under simple ordering prior.

P^* = .900

	$k = 4, \pi = (.1,.6,1,1,1,1)$	$\pi = (.1,.6,1,1,1,1)$	$\pi = (.1,.6,1,1,1,1)$	$\pi = (.1,.6,1,1,1,1)$
	6	7	8	9
PS	.9995	.9995	1.0000	.9995
PI	.0690	.0645	.0545	.0410
PC	.0685	.0640	.0545	.0410
EI	1.4940	1.5345	1.5315	1.7500
EJ	.2120	.2250	.2204	.22940
ES	3.4935	3.5340	3.5315	3.7495
	$k = 5, \pi = (.1,.6,1,1,1,1,2,1)$	$\pi = (.1,.6,1,1,1,1,2,1)$	$\pi = (.1,.6,1,1,1,1,2,1)$	$\pi = (.1,.6,1,1,1,1,2,1)$
	6	7	8	9
PS	.9990	.9995	1.0000	.9995
PI	.0540	.0530	.0295	.0230
PC	.0530	.0520	.0295	.0230
EI	1.9205	1.9700	2.1385	2.4385
EJ	.6140	.6646	.7727	.8577
ES	3.9195	3.9630	4.1385	4.4385
	$k = 4, \pi = (.1,1,1,4,1,1,1,1)$	$\pi = (.1,1,1,4,1,1,1,1)$	$\pi = (.1,1,1,4,1,1,1,1)$	$\pi = (.1,1,1,4,1,1,1,1)$
	6	7	8	9
PS	1.0000	1.0000	1.0000	1.0000
PI	.0375	.0375	.0375	.0375
PC	.0375	.0375	.0375	.0375
EI	.6890	.7475	.8115	.8635
EJ	.6796	.7371	.8031	.8551
ES	.6890	.7475	.8115	.8635
	$k = 5, \pi = (.1,1,1,4,1,1,1,1,2)$	$\pi = (.1,1,1,4,1,1,1,1,2)$	$\pi = (.1,1,1,4,1,1,1,1,2)$	$\pi = (.1,1,1,4,1,1,1,1,2)$
	6	7	8	9
PS	1.0000	1.0000	1.0000	1.0000
PI	.0315	.0315	.0315	.0315
PC	.0315	.0315	.0315	.0315
EI	.9730	.1045	.1360	.1675
EJ	.9635	.1035	.1355	.1670
ES	.9730	.1045	.1360	.1675

CHAPTER II
SAYERS' π^* SELECTION RULES
FOR SELECTING A SUBSET CONTAINING
THE BEST POPULATION

2.1. Introduction

Suppose we have k ($k \geq 2$) independent populations $\pi_1, \pi_2, \dots, \pi_k$. If the random variable X_j associated with π_j has a distribution with known parameters μ_j, σ_j , $j = 1, 2, \dots, k$, first, we give some definitions.

Definition 2.1.1. The population π_1 is the best population of $\pi_1, \pi_2, \dots, \pi_k$ for all $i \neq 1$, if there are more than one population with this condition we arbitrarily tag one of them and call it the best population which is not the best, we called a non-best population.

Assume that we have n_j independent observations $x_{ij}, i = 1, 2, \dots, n_j$, for population π_j , $j = 1, 2, \dots, k$. Let $\hat{\mu}_j$ be the sample mean estimator of μ_j , $j = 1, 2, \dots, k$, so that $\hat{\mu}_j$'s are independent. Let $\hat{\sigma}_j^2$ be the sample variance estimator of σ_j^2 , $j = 1, 2, \dots, k$, for each π_j , we have the following properties of $\hat{\mu}_j$ and $\hat{\sigma}_j^2$:

(a) $\hat{\mu}_j$ is the unbiased estimator of μ_j , $j = 1, 2, \dots, k$.

(b) $\hat{\sigma}_j^2$ is the unbiased estimator of σ_j^2 , $j = 1, 2, \dots, k$.

The general form of the π^* selection rule for selecting a subset of k

$\{1, 2, \dots, k\}$. An action $A \in \mathcal{A}$ is the selection of some subset of the k populations, i.e. $i \in A$ means that x_i is included in the selected subset. An action $A \in \mathcal{A}$ is called a correct selection (CS) if the best population is included in the selected subset A .

Definition 2.1.2. A measurable function γ defined on \mathbb{R}^k is called a selection procedure provided that for each $x \in \mathbb{R}^k$, we have

$$0 < \gamma(x, \delta)$$

and

$$\sum_{A \in \mathcal{A}} \gamma(x, A) = 1$$

where $\gamma(x, A)$ denotes the probability that the subset A is selected when x is observed.

The individual selection probability $\gamma_i(x)$ for the population x_i is then given by

$$\gamma_i(x) = \sum_{A \ni i} \gamma(x, A)$$

where the summation is over all subsets A which contain i . The selection probability $\gamma_i(x)$ takes on only values 0 or 1, i.e., according to the selection procedure $\gamma(x, \delta)$, it completely identifies i . This is what is called a non-randomized procedure.

Definition 2.1.3. Two selection procedures γ and γ' are equivalent if they have the same individual selection probabilities, i.e., if

$$\text{for all } x_i \in \{1, 2, \dots, k\}$$

Hence we can use the following definition, replacing γ by γ' :

Definition 3.1.4. A subset selection rule τ is a measurable function from \mathbb{R}^k to \mathbb{R}^k ,

$$\tau(x) = (\tau_1(x), \dots, \tau_k(x))'$$

with

$$0 \leq \tau_j(x) \leq 1, \quad j = 1, \dots, k.$$

If τ_j 's are 0 or 1, the rule is nonstochastic. Note that, by definition 3.1.1, we have $\sum_{j=1}^k \tau_j(x) \leq 1$.

Suppose our goal is to find a nontrivial subset which contains the best population. A large body of literature exists in the area of subset selection procedures (see Gupta and Sanchezaram (1971), Gupta (1956, 1966) gave maximum-type subset selection procedures. Gupta (1972) studied the performance of Gupta-type maximum procedure, the real-time average procedure (Seal (1956, 1967)) and the Bayesian procedure. Berger (1971c) and Berger and Gupta (1970) proved that Gupta-type maximum procedure is minimax under certain loss functions. In the context of the set-in approach to the subset selection problem, Gold and Puri (1970), Chernoff and Sobay (1971), Buckley and Gray (1972), Gupta and Puri (1972) and Puri (1979) gave different formulation under different loss functions. The loss functions proposed by them are consistent with the set-in approach and the results are applicable to the situation where the elements of the two populations are known, whatever the nature of the loss function may be. In the same litter, Puri and Berger (1972) studied the

in the other hand, an important result by Gupta and Puri (1972) and Puri (1979) is the fact that, if one applies the set-in approach to the selection of a subset of size k from a population of size n , then the effect of the size of the population is negligible.

the different sampling procedures can be derived from the same basic set-up, and we proceed similarly. However, the first two procedures are based on the assumption that all the distributions under consideration have identical parameters, while the last two are based on the assumption that the distributions are of the same type, but only different in their parameters. In the first two cases, the selection problem is independent of the order in which the samples are taken, while in the last two we may wish to relate (modify) the so-called *P*-priorities.

In this chapter we define the posterior-type quantities and introduce two Bayesian selection procedure. $\hat{\pi}_k^B$ and $\hat{\pi}_k^{B*}$ are proposed in section 2.1 and section 2.2 separately, and their properties are discussed in sections 2.3 and 2.4. In section 2.5 we discuss their applications to gamma distributions. In section 2.6, procedure $\hat{\pi}_k^B$ is compared with the well-known $\hat{\pi}_k$ procedure. An application for the problem of selecting the best k elements of the normal distributions $N_{k+1}(\theta_k^*, \sigma^2)$ is given in section 2.7. In section 2.8 we discuss their applications to the selection problems for Poisson distributions and finally we discuss their relation to the selection of gamma distributions. We also compare them with the comparison of the performance of selection procedures by M. and M. (where M_1 and M_2 are the two methods) proposed by a procedure based on single sample and sample mediating, respectively, in Gupta (1956), Gupta and Gaur and Singh (1966). The comparison is done upon Monte Carlo simulation and the results of the comparison are given in terms of the expected expected value and the variance of the expected value of the selection rule, where the individual probabilities of the selection rule are the probabilities of the selection rule being selected. The comparison of the two methods is based on the comparison of the expected value of the selection rule.

2.2. Definitions of the Posterior-P^{*} Condition and the Non-randomized Bayes-P^{*} Procedure

Let $\{1, \dots, k\}$ be the ordered unknown γ_1^{true} . Suppose we have prior distribution p_i for $i \in \{1, \dots, k\}$, then the posterior probability of a correct selection under selection procedure π , given x , is

$$P(\text{CS}_{\pi}, x) = \sum_{i=1}^k p_i(x) p_i$$

where

$$p_i(x) = P(\gamma_i \text{ is the best } | Y = x).$$

It is clear

$$\sum_{i=1}^k p_i(x) \leq 1.$$

Definition 2.2.1. Given a number $P^* (0 < P^* \leq 1)$ and the prior p_i , we say a selection procedure π satisfies the posterior-P^{*} condition if

$$P(\text{CS}_{\pi}, x) \geq P^* \quad \text{for all } x.$$

Remark 2.2.1. The posterior-P^{*} condition is based on the prior distribution p_i and is different from the usual so-called P^{*} condition.

definition 2.2.2. The loss function ψ_{π} is defined by $\psi_{\pi}(x)$, where $|A|$ is the size (number) of populations associated with the selected set A . The loss function ψ_{π} is defined by $\psi_{\pi}(\cdot, A) = \sum_{i \in A} l_i(\cdot)$,

which is the number of the non-best populations selected by π (note). Note that the indicator function

$$\begin{aligned} I_{\{\gamma_1, \dots, \gamma_k\}}(t) &= I_{\{t = \gamma_1\}} + \dots + I_{\{t = \gamma_k\}} \\ &= 0, \text{ otherwise} \end{aligned}$$

definition 2.2.3. Given a number $P^* \frac{1}{k} < P^* \leq 1$ and the prior π , we define the class $\omega_{NR}(\cdot, P^*)$ as follows.

$\omega_{NR}(\cdot, P^*) = \{\cdot\}$ is any non-randomized rule which satisfies the posterior- P^* condition.

For the sake of convenience sometimes we will use ω_{NP} instead of $\omega_{NR}(\cdot, P^*)$.

definition 2.2.4. Given a number $P^* \frac{1}{k} < P^* \leq 1$, a prior π and the loss function L , a selection procedure $\omega_{NP}(\cdot, P^*)$ is called a non-randomized Bayes- P^* procedure (rule) if \cdot is a Bayes rule in the class $\omega_{NP}(\cdot, P^*)$.

Let $p_{(1)}(x) \geq \dots \geq p_{(k)}(x)$ be the ordered $p_i(x)$'s and $\pi_{(i)}$ be the population associated with $p_{(i)}(x)$, $i = 1, \dots, k$, then a subset selection rule \cdot is completely specified by $\omega_{(1)}, \dots, \omega_{(k)}$, where $\omega_{(i)}$ is defined by

$$\omega_{(i)}(x) = P\{x \in S_i\}, \quad i = 1, \dots, k.$$

Next, we propose a non-randomized selection rule which belongs to

$\omega_{NP}(\cdot, P^*)$.

definition 2.2.5. Given a number $P^* \frac{1}{k} < P^* \leq 1$, a prior π and a prior distribution π , the selection rule ω_{NP} is defined by $\omega_{NP}(x_1, x_2, \dots, x_k)$, where

$$\omega_{NP}(x_1, x_2, \dots, x_k) = \begin{cases} 1, & \text{if } x_i = \hat{x}(x) \\ \frac{1}{k}, & \text{if } x_i = \bar{x}(x) \\ 0, & \text{otherwise} \end{cases}$$

and $j(x)$ is the maximum integer such that

$$\sum_{i=1}^{j-1} p_{\{i\}}(x) < p^*.$$

Lemma 2.2.1. $\frac{B}{NR} \in \mathcal{L}_{NR}$.

Proof. Follows from the definition of $\frac{B}{NR}$.

Theorem 2.2.1. Given a number $p^* \in (p^*, 1)$, the prior π and the function L_1 , the selection procedure $\frac{B}{NR}$ is a non-randomized Bayes rule.

Proof. It is sufficient to show that the selection procedure $\frac{B}{NR}$ has the smallest posterior risk in the class $\mathcal{L}_{NR}(p^*)$. Given the observation $X = x$. Let the posterior risk of $\pi \in \mathcal{L}_{NR}(p^*)$ be $r(x, \pi)$,

$$r(x, \frac{B}{NR}) = k - j + 1$$

and

$$\sum_{i=j+1}^k p_{\{i\}}(x) < p^*$$

for some j , $1 \leq j \leq k$.

Hence the inequality

$$r(x, \pi) \geq r(x, \frac{B}{NR})$$

is not true for any $\pi \in \mathcal{L}_{NR}(p^*)$, $\pi \neq \frac{B}{NR}$. Therefore, the result follows.

Theorem 2.2.2. Theorem 2.2.1 also holds when we replace the loss L_1 by

Proof. Under the loss function L_2 , the posterior risk of the selection procedure $\pi \in \mathcal{L}_{NR}^*(\pi, p^*)$ is

$$r(x, \pi) = \sum_{i=1}^k \pi(i)(x)[i - p_i(x)], \text{ given } x \in \mathcal{X}.$$

By Theorem 2.2.1, we have

$$\sum_{i=1}^k \pi_{NR(i)}^B(x) + \sum_{i=1}^k \pi(i)(x) = 1.$$

If

$$\sum_{i=1}^k \pi_{NR(i)}^B(x) < \sum_{i=1}^k \pi(i)(x)$$

then by definition of π_{NR}^B , we have

$$\sum_{i=1}^k \pi_{NR(i)}^B(x)p_i(x) < \sum_{i=1}^k \pi(i)(x)p_i(x).$$

On the other hand, if

$$\sum_{i=1}^k \pi_{NR(i)}^B(x) > \sum_{i=1}^k \pi(i)(x)$$

then

$$\begin{aligned} \sum_{i=1}^k \pi_{NR(i)}^B(x) + \sum_{i=1}^k \pi(i)(x) &= 1 \\ &> \sum_{i=1}^k \pi(i)(x)(1 - p_i(x)). \end{aligned}$$

Therefore, we have

$$r(x, \pi_{NR}^B) < r(x, \pi) \text{ for all } \pi \in \mathcal{L}_{NR}^*(\pi, p^*).$$

Corollary 2.2.1. For a given prior π and the loss function

$L = t_1 \mathbb{1}_1 + t_2 \mathbb{1}_2$ where $t_1, t_2 > 0$, then π_{NR}^B is a non-dominated Bayes

rule wrt the loss function ℓ , for all x_1, \dots, x_k .

Proof. For the given prior π and the loss function ℓ , the Bayes risk of any procedure π^* on \mathcal{X}_{BFR} is, given $x = (x_1, \dots, x_k)$,

$$\begin{aligned} R(x, \pi^*) &= C_1 \sum_{i=1}^k \ell_{\pi^*}(x_i) + C_2 \sum_{i=1}^k \ell_{\pi^*}(x_i) \mathbb{P}_{\pi^*}(\{x_i\}) \\ &= C_1 \sum_{i=1}^k p_{\pi^*}(x_i) \ell_{\pi^*}(x_i) + C_2 \sum_{i=1}^k \frac{p_{\pi^*}(x_i)}{\mathbb{P}_{\pi^*}(\{x_i\})} \ell_{\pi^*}(x_i) \\ &= \ell_{\pi^*}(x, \pi_{\text{BFR}}) \end{aligned}$$

wrt the loss function ℓ .

Hence π_{BFR}^* is a Bayes-E * -rule wrt the loss function ℓ for all x .

3.3. Proposed Bayes-E * Procedure π_{BEP}^* in general

Suppose we are interested in the randomized subset selection rule, and we would like to find such a rule which also satisfies the E * -condition and has the minimum risk wrt the loss function ℓ , given the prior distribution π .

Definition 3.3.1. Given a prior π , we define a Bayes-E * -subset selection rule π^* , in which all rules satisfy the uniform E * -condition for any given observation $x = (x_1, \dots, x_k)$, that is,

$$\ell_{\pi^*}(x, \pi^*) = \inf_{\pi \in \Pi} \ell_{\pi^*}(x, \pi) \quad \text{for all } x.$$

Definition 3.3.2. Given a number $C_1 \geq 0$, a prior π and a loss function ℓ , a selection procedure π_{BEP}^* is called a Bayes-E * -rule if the procedure π_{BEP}^* is a Bayes rule to the class $\mathcal{X}_{\text{BEP}}^*$.

For the sake of convenience, sometimes we will use π^* instead of $\pi(\cdot, p^*)$.

Definition 2.3.2. We define a subset selection procedure π^B as follows.

Given a prior π and observation $x = x_1, \dots, x_B$ is defined by

$$\pi^B(1), \pi^B(2), \dots, \pi^B(k),$$

where

$$\pi^B(k)(x) = 1$$

and

$$\begin{aligned} \pi^B(j)(x) &= 1, \quad \text{if } \sum_{i=j+1}^k p_{[i]}(x) < p^*, \quad j \neq k \\ &= 0, \quad \text{if } \left\{ \begin{array}{l} \sum_{i=j+1}^k p_{[i]}(x) > p^* \\ \sum_{i=j+1}^k p_{[i]}(x) + \frac{k-j}{k-j+1} p_{[j]}(x) > p^* \end{array} \right. \\ &= 0, \quad \text{otherwise.} \end{aligned}$$

Example. If $k = 5$, $p^* = .90$ and the posterior probabilities are $p_{[1]}(x) = .30$, $p_{[2]}(x) = .15$, then we select the population $[1]$ with probability $p_{[1]}(x)$ with probability $.15$. And we select $[2]$ with probability $.15$, where $p_{[2]}(x)$ is given by

$$p_{[2]}(x) = \min\left(1, \frac{.30}{.15}\right) = 2.00$$

$$\frac{1}{1 - \frac{.30}{.15}} = \frac{1}{1 - 2.00} = \frac{1}{-1.00} = -1.00$$

By Definition 2.3.2 we have

$$\min_{\pi \in \Pi} B_{\pi}(x, \cdot, \cdot) = \max_{i=1}^k \pi_{(i)}(\cdot) \pi_{(i)}(\cdot) = \pi^{*} \text{ at } \pi_{(k)}(\cdot)$$

hence we have the following Lemma.

Lemma 2.3.3. $\pi^{*} \in \arg \max_{\pi \in \Pi} B_{\pi}(x, \cdot, \cdot)$.

DEFINITION 2.3.4. we define a subclass $\omega^*(\pi, P^*)$ of class $\omega(\pi, P)$ by

$$\omega^*(\pi, P^*) = \{\pi \in \omega(\pi, P) : \pi_{(1)}(x) \leq \pi_{(2)}(x) \leq \dots \leq \pi_{(k)}(x)\}$$

where $\pi_{(1)}(x) \leq \dots \leq \pi_{(k)}(x)$ are the ordered $\pi_i(x)$'s.

By the definition of $\omega^*(\pi, P^*)$ we have the following Lemma.

Lemma 2.3.5. $\pi^{*}, \beta, \gamma, \gamma_{(1)}^k \in \omega^*(\pi, P^*)$.

Lemma 2.3.6. for all $\pi \in \omega(\pi, P^*)$ there exists $\beta \in \omega^*(\pi, P^*)$ such that $\beta(x, \cdot) \leq \pi(x, \cdot)$ wrt the loss function L_P , for all x .

Theorem 2.3.7. Selection procedure of a Bayesian procedure in $\omega(\pi, P^*)$ wrt the loss function L_P .

Proof. Given the observation $x = x_k$ and any selection procedure

$\omega(\pi, P^*) \ni \pi_{(k)}(x) \leq \pi_{(k+1)}(x) \leq \dots \leq \pi_{(n)}(x)$ and $\pi_{(1)}(x) \leq \dots \leq \pi_{(k-1)}(x)$ if $k \neq K$, hence we have

$$\min_{\pi \in \Pi} B_{\pi}(x, \cdot, \cdot) = \pi_{(k)}(x), \text{ for all } \pi \in \omega(\pi, P^*)$$

If $\pi_{(k)}(x) = \pi^{*}$, then we have $\pi_{(k)}(x) \leq \pi_{(k+1)}(x) \leq \dots \leq \pi_{(n)}(x)$ and $\pi_{(1)}(x) \leq \dots \leq \pi_{(k-1)}(x)$.

We will now show that for any π ,

$$\pi_{(k)}(x) \leq \pi_{(k)}(x_k)$$

implies

$$P(\text{CS}_{i+1}^A, x) \geq P(\text{CS}_{i+1}^B, x) = \pi_{i+1}^{*},$$

that is,

$$\sum_{j=1}^k \pi_{i+1,j}^A(x) \geq \sum_{j=1}^k \pi_{i+1,j}^B(x)$$

implies

$$\sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^A(x) \geq \sum_{j=1}^k \pi_{i+1}^B(j) \rho_{i+1,j}^B(x).$$

For any $0 < t < \sum_{j=1}^k \pi_{i+1,j}^B(x)$, we have $t = a + b$ where a is a positive integer and $0 < b < 1$.

It is easy to see that the maximum posterior probability of a non-optimal selection of procedure j with $\sum_{j=1}^k \pi_{i+1,j}^A(x) = c$ is

$$\sum_{j=k-a+1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^A(x) + \pi_{i+1,k-a}^A(x),$$

and it is less than $\sum_{j=1}^k \pi_{i+1,j}^B(x) \rho_{i+1,j}^B(x)$, since $c < \sum_{j=1}^k \pi_{i+1,j}^B(x)$. Therefore π_{i+1}^A is Bayes-P* procedure in $x \in \mathcal{X}^A$.

Lemma A. Given the loss function L_{α} for all $\alpha \in \mathcal{A}$, there exists a Bayes-P* such that it is better than any other Bayes-P* procedure in the posterior risk with the prior and the loss function.

Proof. Given π_{i+1}^A , let $\pi_{i+1}^B = \pi_{i+1}^A - \pi_{i+1}^A$. Then

$$\begin{aligned} \sum_{j=1}^k \pi_{i+1}^B(j) \rho_{i+1,j}^B(x) &= \sum_{j=1}^k (\pi_{i+1}^A - \pi_{i+1}^A)(j) \rho_{i+1,j}^B(x) \\ &= \sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^B(x) - \sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^B(x) \\ &= \sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^A(x) - \sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^B(x) \\ &= \sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^A(x) - \sum_{j=1}^k \pi_{i+1}^A(j) \rho_{i+1,j}^A(x) \\ &= 0. \end{aligned}$$

$$\sum_{i=1}^k \pi_i^{(x)} v_i^{(x)}$$

$$= p^*$$

hence $\pi^* \in \omega^*(\cdot, p^*)$.

Now,

$$\begin{aligned} \pi(s_1, \cdot) &= \sum_{i=1}^k \pi_i(s_1)(1 - \pi_{i+1}(s_1)) \\ &= \sum_{i=1}^k \pi_i(s)(1 - \pi_{i+1}(s)) \\ &\leq \sum_{i=1}^k \pi_i(s)(1 - \pi_{i+1}^*(s)) \\ &= \pi(s_1, \cdot). \end{aligned}$$

Hence the proof is complete.

Theorem 2.3.2. Given the prior π and the observation s_1 , the procedure π^* is a Bayes- p^* procedure in the class $\omega^*(\cdot, p^*)$ when the prior distribution is π_1 .

Proof. By Lemma 2.2.4, it is sufficient to show that

$$\pi(s_1, \frac{d}{ds}) = \min_{\pi' \in \omega^*(\cdot, p^*)} \pi(s_1, \frac{d}{ds})$$

where

$$\begin{aligned} \pi(s_1, \frac{d}{ds}) &= \sum_{i=1}^k \pi_i(s_1)(1 - \pi_{i+1}(s_1)) \\ &= \sum_{i=1}^k \pi_i(s) \frac{\pi_i(s)}{\pi_i(s) + \pi_{i+1}(s)} (1 - \pi_{i+1}^*(s)). \end{aligned}$$

□

$$i_0 = \min_{1 \leq i \leq k} \left(\frac{B}{i} \right)^{\lambda^k} - 1$$

$$\hat{A}_1 = \left(\frac{B}{i_0} \right)^{\lambda^k} \left(x \right) + \left(\frac{B}{i_0} \right)^{\lambda^k} \left(x \right) > 0$$

$$\hat{A}_{\mu} = \left(\frac{B}{i_\mu} \right)^{\lambda^k} \left(x \right) + \left(\frac{B}{i_\mu} \right)^{\lambda^k} \left(x \right) > 0, \quad \forall \mu,$$

then $A_1 + i_0 \geq 1$.

And we have

$$d_1 = \max \left(\hat{A}_1 - i_0, -1 \right), \quad \text{if } \left(\frac{B}{i_0} \right)^{\lambda^k} \leq \left(\frac{B}{i_1} \right)^{\lambda^k},$$

$$d_0 = -1, \quad \text{if } \left(\frac{B}{i_0} \right)^{\lambda^k} > \left(\frac{B}{i_1} \right)^{\lambda^k}$$

$$d_2 = \min \left(\hat{A}_2 - i_0, -1 \right), \quad \text{if } \left(\frac{B}{i_0} \right)^{\lambda^k} \leq \left(\frac{B}{i_2} \right)^{\lambda^k},$$

$$i_0 + 1, \quad \text{if } \left(\frac{B}{i_0} \right)^{\lambda^k} > \left(\frac{B}{i_2} \right)^{\lambda^k}$$

hence $d_1 = d_2$.

Therefore, we have

$$\begin{aligned} e(x, v^k) - e(x, v^{\bar{k}}) &= \sum_{i=1}^k \left(\frac{B}{i} \right)^{\lambda^k} \left(x \right) + \left(\frac{B}{i} \right)^{\lambda^k} \left(x \right) - \left(\frac{B}{i} \right)^{\lambda^{\bar{k}}} \left(x \right) \\ &\quad - \sum_{i=k+1}^{\bar{k}} \left(\frac{B}{i} \right)^{\lambda^k} \left(x \right) - \left(\frac{B}{i} \right)^{\lambda^k} \left(x \right) + \left(\frac{B}{i} \right)^{\lambda^{\bar{k}}} \left(x \right) \\ &= \sum_{i=k+1}^{\bar{k}} \left(\frac{B}{i} \right)^{\lambda^k} \left(x \right) - \left(\frac{B}{i} \right)^{\lambda^{\bar{k}}} \left(x \right) \\ &= \sum_{i=k+1}^{\bar{k}} \left(\frac{B}{i} \right)^{\lambda^k} \left(x \right) - \left(\frac{B}{i} \right)^{\lambda^{\bar{k}}} \left(x \right) \end{aligned}$$

$$\sum_{i \in A_1} \frac{c_i^B(x)}{c_i^A(x)} \frac{\frac{B}{A}(x) - c_i^A(x)}{\frac{B}{A}(x) + c_i^A(x)} = \frac{1}{\sum_{i \in A_1} \frac{c_i^B(x)}{c_i^A(x)}}.$$

$$\star(1 + \frac{B}{A}(x))$$

$$\sum_{i \in A_1} \frac{c_i^B(x)}{c_i^A(x)} + \frac{B}{A}(x)(1 + \frac{B}{A}(x))$$

as required by Theorem 2.3.1.

Corollary 2.3.4. Procedure \star is a Bayes- c^* rule in L^A if and only if the loss function $L = c_1^A + c_2^B$, $c_1, c_2 > 0$.

Proof. Similar to Corollary 2.2.1, hence omitted.

2.4. Properties of $\frac{B}{A}$ and $\frac{B}{A^2}$

In this section we discuss some properties of selection functions $\frac{B}{A}$ and $\frac{B}{A^2}$. The following definition of the ordering of distributions was introduced by Lehmann (1953) and further discussed by Lehmann and Bain (1975).

Definition 2.4.1. A subset $A \subseteq \mathbb{R}^k$ is a set of distributions if $c_i > 0$ and $c_i > c_j$ for all $i \neq j$, $i, j \in A$, otherwise no.

Definition 2.4.2. A family of distributions A is called a $\frac{B}{A}$ -family if there exists a function $\frac{B}{A}(x)$ such that the following two conditions hold for all $x \in \mathbb{R}^k$ and $a_1, a_2 \in A$ (Lehmann and Bain 1975).

$$\frac{B}{A}(x) = \frac{\int_A dP(x, a)}{\int_A dP(a)}$$

for all x in the domain of $\frac{B}{A}$.

Let $f(\cdot, \gamma_j)$ be the p.d.f. of population γ_j . Let π_{ij} be the prior where γ_j 's are mutually independent. Suppose for $x = x_i$, we have absolutely continuous posterior c.d.f. $q_i(\cdot | x)$. Hence we can write the p.d.f. as

$$q(x) = \prod_{j=1}^k q_j(\gamma_j | x) = \prod_{j=1}^k q_j(\gamma_j | x_i),$$

Let $a_i(\cdot, x_i)$ be the posterior c.d.f. associated with $\gamma_i + \gamma_{i+1}, i = 1, \dots, k$.

Definition 7.4.3. The absolutely continuous posterior c.d.f., $a_i(\cdot, x_i), i = 1, \dots, k$, have the generalized (strictly) stochastic increasing property (G-SIP) if for any $i, j, l, i < j < l < k$, $a_i(\cdot, x_i) \leq$

$$a_l(\cdot, x_j) \quad \text{and} \quad a_j(\cdot, x_i).$$

Note that if $a_i(\cdot, x_i) \leq a_j(\cdot, x_j) \leq a_l(\cdot, x_l)$ for all $i < j < l$, then the SIP is the usual SIP.

Definition 7.4.4. A selection procedure γ is monotone decreasing if and only if for every $x \in \mathbb{R}^k$, $x_i > x_j$ implies $\gamma_i < \gamma_j$. In other words, γ is monotone with the increasing order of x almost surely with probability zero.

Theorem 7.4.1. If the prior π is such that we have $a_i(\cdot, x_i), i = 1, \dots, k$ are independent posterior c.d.f. and $a_i(\cdot, x_i) \leq a_j(\cdot, x_j)$ for all $i < j$, then for every $x \in \mathbb{R}^k$, $x_i > x_j$ implies

$$\gamma_i < \gamma_j \quad \text{almost surely.}$$

Since both the selection procedure γ and $a_i(\cdot, x_i)$ are absolutely continuous, the selection procedure γ is monotone decreasing.

Proof.

$$\begin{aligned}
 p_i(x) &= P(\tau_{ij} < \tau_{ik} | x) = \int_{\mathbb{R}^d} \frac{\partial}{\partial t} P(\tau_{ij} < t | x) d\mu_j(t) d\mu_k(t) \\
 &= \int_{\mathbb{R}^d} \int_{\mathbb{R}^d} \frac{\partial}{\partial t} P(\tau_{ij} < t | x) M_j(t|x) d\mu_j(t) d\mu_k(t) \\
 &= \int_{\mathbb{R}^d} \int_{\mathbb{R}^d} \frac{\partial}{\partial t} P(\tau_{ij} < t | x) M_j(t|x) \mu_j(t) d\mu_k(t) \\
 &= \int_{\mathbb{R}^d} \frac{\partial}{\partial t} P(\tau_{ij} < t | x) \mu_j(t) \mu_k(t) dt \\
 &= \int_{\mathbb{R}^d} \frac{\partial}{\partial t} [P(\tau_{ij} < t | x) \mu_j(t)] \mu_k(t) dt \\
 &= \int_{\mathbb{R}^d} \frac{\partial}{\partial t} P(\tau_{ij} < t | x) \mu_j(t) \mu_k(t) dt
 \end{aligned}$$

Since

$$B_j(x) = \frac{b}{c_j}(x) - \gamma + \delta - \mu_j(x) - \mu_k(x)$$

and

$$\frac{B}{\text{NR}_j}(x) = \frac{B}{c_j}(x) - \gamma + \delta - \mu_j(x) - \mu_k(x)$$

Therefore, the procedures $\frac{P}{\cdot}$ and $\frac{B}{\text{NR}_j}$ are ordered to

Under (VII) assumptions, we can relate the population parameter $c_j = c_k = c$ by hence we have $\mu_j(x) = \mu_k(x)$ and $\frac{c_j}{c_k} = \frac{c_j}{c_k} = 1$.
 Tsiatis (1970) defined the "last" procedure to be the one that is more definitive (VII). A set of rules for ordering the last procedures are:
 (i) if all x such that $\frac{B}{c_j}(x) > \frac{B}{c_k}(x)$ are contained in \mathcal{C}_{jk} then $\frac{B}{c_j} > \frac{B}{c_k}$
 (ii) if $\mathcal{C}_{jk} = \emptyset$

we call $\frac{B}{c_j}(x)$ the "last" value to everywhere else in \mathcal{C}_{jk} and the corresponding $\frac{B}{c_k}(x)$ is called the "second last" value. In our procedure, we

Definition 7.4.6. A selection rule γ_j is called translation invariant if for all $x \in \mathbb{R}^k$, and for all $c \in \mathbb{R}$

$$\gamma_j(x_1 + c, \dots, x_k + c) = \gamma_j(x_1, \dots, x_k) \quad \text{if } x_1, \dots, x_k \neq 0$$

Definition 7.4.7. A selection rule γ_j is called scale-invariant if for all $x \in \mathbb{R}^k$, and for all $c > 0$

$$\gamma_j(cx_1, \dots, cx_k) = \gamma_j(x_1, \dots, x_k) \quad \text{if } x_1, \dots, x_k \neq 0$$

Theorem 7.4.2. If the posterior distributions $\alpha_{j+1}, \dots, \alpha_n$ (for $j+1, \dots, n$) have the shift property, then both selection procedures γ_j and $\gamma_{\bar{j}}$ are "justifiable".

PROOF. It is sufficient to show that

$$D_j(x) = D_{j+1}(x) \quad \text{whenever } x_j^T = x_{j+1}^T$$

$$(0 < x_1 < x_2 < \dots < x_j < x_{j+1} < \dots < x_n)$$

or equivalently,

$$\gamma_j(x_1, \dots, x_{j+1}) = \gamma_{j+1}(x_1, \dots, x_{j+1})$$

$$\gamma_{j+1}(x_1, \dots, x_{j+1}) = \gamma_{j+2}(x_1, \dots, x_{j+2})$$

$$\vdots$$

$$\gamma_{n-1}(x_1, \dots, x_{n-1}) = \gamma_n(x_1, \dots, x_n)$$

$$\gamma_n(x_1, \dots, x_n) = \gamma_{\bar{j}}(x_1, \dots, x_n)$$

$$x = (x_1, \dots, x_n)$$

Let us consider the following two cases:

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Definition 2.4.8. Given a number $P^* \left(\frac{1}{k} < P^* < 1 \right)$, $X = x$ and a prior τ ; for any selection procedure $\psi \in \Delta(\tau, P^*)$ the ratio of the posterior probability $P(CS|\psi, x)$ and the posterior expected selected size $E(S|\psi, x)$ is called the posterior-efficiency of ψ and is denoted by $EFF(\psi|x)$.

$$EFF(\psi|x) = \frac{P(CS|\psi, x)}{E(S|\psi, x)}.$$

If $EFF(\psi|x) \geq EFF(\psi'|x)$ for all $\psi' \in \Delta$ and all x , then the selection procedure ψ is called "posterior most efficient" (PME) selection procedure in $\Delta(\tau, P^*)$.

Theorem 2.4.3. The non-randomized posterior- P^* selection procedure ψ_{NR}^B is the PME selection procedure in $\Delta_{NR}^I(\tau, P^*) = \Delta_{NR}$, given τ, P^* .

Proof. By Lemma 2.2.1, for all $\psi \in \Delta_{NR}$

$$\exists \psi' \in \Delta_{NR}^I \Rightarrow EFF(\psi'|x) \geq EFF(\psi|x) \quad \forall x,$$

hence it is sufficient to show that:

Given $\tau(\underline{x}), P^*, \underline{x}$, $EFF(\psi_{NR}^B|x) \geq EFF(\psi'|x)$ for all $\psi' \in \Delta_{NR}^I(\tau, P^*)$.

We know that, in $\Delta_{NR}(\tau, P^*)$ hence in $\Delta_{NR}^I(\tau, P^*)$, ψ_{NR}^B always has minimum selected size, i.e. $\forall x, \sum_{i=1}^k \psi_{NR}^B(x) + c = \sum_{i=1}^k \psi_i^I(x)$ for some integer c , $0 \leq c \leq k - 1$.

$$\begin{aligned} EFF(\psi'|x) &= \frac{\sum_{i=1}^k \psi_i^I(x) p_{[i]}(x)}{\sum_{i=1}^k \psi_i^I(x)} \\ &< \frac{\sum_{i=1}^k \psi_{NR}^B(x) p_{[i]}(x) + p_{[k-s-c+1]}(x) + \dots + p_{[k-s]}(x)}{\sum_{i=1}^k \psi_i^B + c} \end{aligned}$$

if $\psi_{NR}^B(x) = (0, \dots, \underbrace{0}_{s \text{ terms}}, 1, \dots, 1)$.

$$\begin{aligned} \text{EFF}(\psi' | \underline{x}) &\leq \frac{\sum_{i=k-s+1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x}) + c p_{[k-s]}(\underline{x})}{\sum_{i=1}^k \psi_{NR(i)}^B(\underline{x}) + c} \\ &\leq \frac{\sum_{i=1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x})}{\sum_{i=1}^k \psi_{NR(i)}^B(\underline{x})} \\ &= \text{EFF}(\psi_{NR}^B | \underline{x}). \end{aligned}$$

The last inequality is obtained by

$$\begin{aligned} \sum_{i=1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x}) &= \sum_{i=k-s+1}^k \psi_{NR(i)}^B(\underline{x}) p_{[i]}(\underline{x}) \\ &\geq \left(\sum_{i=k-s+1}^k \psi_{NR(i)}^B(\underline{x}) \right) p_{[k-s]}(\underline{x}). \end{aligned}$$

Theorem 2.4.4. The randomized selection procedure ψ^B is the PME procedure in $\mathcal{A}(\tau, P^*) = \mathcal{A}$ for given τ, P^* .

Proof. It suffices to show that, given $\tau, P^*, \underline{x} = \underline{x}$,

$$\text{EFF}(\psi^B | \underline{x}) \geq \text{EFF}(\psi' | \underline{x}), \quad \forall \psi' \in \mathcal{A}'.$$

Suppose $\psi^B(\underline{x}) = (0, \dots, \underbrace{v, 1, \dots, 1}_{s \text{ terms}}) \quad 0 \leq v < 1, 1 \leq s \leq k - 1$.

By theorem 2.3.1 there exists $c > 0$ such that

$$\sum_{i=1}^k \psi(i)(\underline{x}) + c = \sum_{i=1}^k \psi'(i)(\underline{x}).$$

If $0 \leq c < 1$, then

$$\begin{aligned} \text{EFF}(\psi' | x) &= \frac{\sum_{i=1}^k \psi'_i(x) p_{[i]}(x)}{\sum_{i=1}^k \psi'_i(x)} \\ &\leq \frac{\sum_{i=1}^k \psi_i^B(x) p_{[i]}(x) + cp_{[k-s]}(x)}{\sum_{i=1}^k \psi_i^B(x) + c} \\ &\leq \frac{\sum_{i=1}^k \psi_i^B(x) p_{[i]}(x)}{\sum_{i=1}^k \psi_i^B(x)} \\ &= \text{EFF}(\psi^B | x). \end{aligned}$$

If $1 \leq c = v' + t + (1-v)$, $t > 0$ integer, $0 < v' < 1$ then

$$\begin{aligned} \sum_{i=1}^k \psi'_i(x) p_{[i]}(x) &= \sum_{i=k-s+1}^k \psi_i^B(x) p_{[i]}(x) + v' p_{[k-s-t+1]}(x) \\ &\quad + p_{[k-s-t]}(x) + \dots + (1-v) p_{[k-s]}(x) \\ &\geq \sum_{i=1}^k \psi_i^B(x) p_{[i]}(x) + cp_{[k-s]}(x) \end{aligned}$$

hence by the same argument as above we have

$$\text{EFF}(\psi' | x) \geq \text{EFF}(\psi^B | x).$$

Since x is arbitrary, the result holds for all x .

2.5. Applications to Normal Model

Suppose we have k populations π_1, \dots, π_k ; population π_i has distribution $N(\mu_i, \sigma_i^2)$, where σ_i 's are known and μ_i 's are unknown. Assume that we

have independent observations x_{i1}, \dots, x_{in_i} , $i = 1, \dots, k$. Let

$$\bar{x}_i = \frac{1}{n_i} \sum_{j=1}^{n_i} x_{ij} \text{ and let } \underline{x} = (\underline{x}_1, \dots, \underline{x}_k).$$

Suppose we are interested in selecting a subset containing the best (the population having the largest mean) under the posterior-P* condition, wrt some prior $\pi = \pi(\mu)$. Then to find a Bayes-P* selection procedure is equivalent, in some sense, to finding $p_i(\underline{x})$, which is the posterior probability of the event $\{\mu_i \text{ is the best}\}$, given observations $\underline{x} = \underline{x}$, wrt a given prior π , for all $i = 1, \dots, k$.

Case I. Assume that we have a common sample size n and a common known variance σ^2 .

Ia. Suppose we have no prior information about the unknown parameters, and use the "non-informative" (Box and Tiao (1973)) or "locally uniform" prior $p(\mu_i) \propto c$ for each population.

The posterior density function g_i of μ_i , given \underline{x} is the normal density with mean \bar{x}_i and variance σ^2/n , i.e.,

$$g_i(\mu_i | \underline{x}) = \frac{c}{\sqrt{2\pi}} \exp \left(-\frac{n(\mu_i - \bar{x}_i)^2}{2\sigma^2} \right).$$

Hence

$$\begin{aligned} p_{[i]}(\underline{x}) &= P(\mu_{(i)} = \mu_{[k]} | \underline{x} = \underline{x}) \\ &= \int_{-\infty}^{\mu_{[i]}} \frac{1}{\sqrt{2\pi}} \exp \left(-\frac{n}{2\sigma^2} (t + \frac{\bar{x}_i}{\sigma})^2 \right) dt \\ &\quad i = 1, \dots, k. \end{aligned}$$

Here $\mu_{(i)}$ is the quantity corresponding to the i^{th} largest observation $x_{[i]}$.

Ib. If μ_i 's are independent and have the identical prior distribution $N(\theta_0, \sigma_0^2)$ and $x_i | \mu_i \sim N(\mu_i, \sigma_1^2/n)$, then it is well known that the posterior density function g_i of μ_i , given $\underline{x} = \underline{x}$ is

$$g_i(\mu_i | \underline{x}) \sim N(\bar{\theta}_{x_i}, \xi^2) \text{ with SIP property}$$

where

$$\bar{\theta}_{x_i} = \xi^2 (\sigma_0^{-2} \theta_0 + n \sigma_1^{-2} x_i)$$

$$\xi^2 = (\sigma_0^{-2} + n \sigma_1^{-2})^{-1}.$$

Hence

$$p_{[i]}(\underline{x}) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi(t + \xi n \sigma_1^{-2} (x_{[i]} - x_{[j]})) d\phi(t).$$

The last expression for $p_{[i]}(\underline{x})$ is the same as that for the non-informative prior whenever $\sigma_0 \rightarrow \infty$.

Since $p_{[i]}(\underline{x}) = p_{[i]}(\underline{x} + b)$ and since the normal distribution has the strictly SIP, it follows that ψ^B and ψ_{NR}^B are "just" a.e. and translation-invariant in both case Ia and Ib.

Case II. Variance σ_i 's are known but σ_i 's and n_i 's are not all equal.

IIa. Using the non-informative prior $p(\mu_i) = c$, $i=1, \dots, k$, we have

$$p_{(i)}(\underline{x}) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi(t + \frac{v_{(i)}}{v_{(j)}} + \frac{x_{[i]} - x_{[j]}}{v_{(j)}}) d\phi(t)$$

where $v_{(i)} = \frac{\sigma_{(i)}}{n_{(i)}}$, $i = 1, \dots, k$. $p_{(i)}$, $v_{(i)}$ and $n_{(i)}$ are corresponding to $x_{[i]}$ and we have the following theorem.

Theorem 2.5.1. $p_{(i)}(x)$ is non-decreasing in i , i.e., $p_{(i)}(x) = p_{[i]}(x)$.

Remark 2.5.1. From the above formula of $p_{(i)}(x)$, it is easy to see, increasing the sample size of the non-best populations will increase the probability that the best population to be selected, however, before doing this, we don't know which one is the best one.

In this case ψ^B and ψ_{NR}^B are "just" a.e. and translation-invariant.

Case III. Assume that priors are independent but not identical normal distributions, namely, $\mu_i \sim N(\theta_i, \sigma_{0i}^2)$, where θ_i 's are not all equal; if the conditional distribution of X_i , given μ_i , is $N(\mu_i, \frac{\sigma_{1i}^2}{n_i})$, then the posterior density of μ_i , given $X_i = x_i$ is $g_i(\mu_i | x_i)$, which is the probability density function of normal distribution $N(\bar{\theta}_{x_i}, \varepsilon_i^2)$ where

$$\bar{\theta}_{x_i} = \varepsilon_i^2 (\sigma_{0i}^{-2} \theta_i + n_i \sigma_{1i}^{-2} x_i)$$

$$\varepsilon_i^2 = (\sigma_{0i}^{-2} + \sigma_{1i}^{-2} n_i)^{-1}.$$

Hence we have

$$p_i(x) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi\left[t - \frac{\varepsilon_j}{\varepsilon_i} + \frac{1}{\varepsilon_j} (\bar{\theta}_{x_i} - \bar{\theta}_{x_j})\right] d\phi(t).$$

If $\sigma_{0i} = \sigma_0$, $\sigma_{1i} = \sigma_1$ and $n_i = n$, $i = 1, \dots, k$, then

$$\varepsilon_i = \varepsilon = (\sigma_0^{-2} + \sigma_1^{-2} n)^{-1} \quad i = 1, \dots, k$$

and

$$p_i(x) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi\left[t + \left(\frac{\theta_i - \theta_j}{\sigma_0^2} + \frac{n(x_i - x_j)}{\sigma_1^2}\right)\right] d\phi(t).$$

Case IV. The General Normal Model

Here we consider a more general prior. Suppose we have k populations, common sample size n for each population, and common known variance $\sigma^2 > 0$. The observation can reduce to $\underline{X} = (X_1, \dots, X_k)$ where

$$X_i = \frac{1}{n} \sum_{j=1}^n X_{ij}, \text{ by sufficiency.}$$

The "Normal Model" is defined as follows:

$$(a) \quad X_i | \underline{\mu} \sim N(\underline{\mu}, qI), \quad q = \frac{\sigma^2}{n}$$

where I is the $k \times k$ identity matrix.

So the X 's are (conditionally) independent when $\underline{\mu}$ is given.

$$(b) \quad \underline{\mu} \sim N(\theta_0 \underline{1}, \gamma I + tU)$$

where $\theta_0 \in \mathbb{R}$, $\gamma > 0$, $t > -\frac{\gamma}{k}$,

$$\underline{1} = (1, \dots, 1) \quad \text{and} \quad U = \underline{1}' \underline{1}.$$

Here $\gamma > 0$ and $t > -\frac{\gamma}{k}$ are necessary and sufficient for $\gamma I + tU$ to be positive definite. This model was chosen by Chernoff and Yahav (1977) ($t > 0$), Gupta and Hsu (1978) and Miescke (1979).

By (a) and (b) we get the posterior distribution of $\underline{\mu}$, given $\underline{X} = \underline{x}$, and the distribution of $\underline{\lambda}$ as follows:

$$\underline{\mu}' \underline{x} \sim N(q, aI + bU)$$

where

$$q = \gamma(q+\gamma)^{-1} \underline{x} + qt((q+\gamma)(q+\gamma+kt))^{-1} \underline{x}' U + q(q+\gamma+kt)^{-1} \underline{1}$$

$$a = \gamma q(q+r)^{-1}$$

$$b = q^2 t(q+\gamma)^{-1} (q+\gamma+kt)^{-1}$$

$$\underline{x} \sim N(\underline{m}, (q+\gamma)I + tU)$$

Lemma 2.5.1. Let $\underline{Y} \sim N(\underline{\mu} + \rho \underline{1}, aI + bU)$ with $\underline{\mu} \in \mathbb{R}^k$, $\rho \in \mathbb{R}$, $a > 0$ and $b > -a/k$. Then there exists a random vector $\underline{Z} \sim N(\underline{\mu}, aI)$ such that $h(Y) = h(Z)$ everywhere for every translation-invariant $h: \mathbb{R}^k \rightarrow \mathbb{R}^k$.

Proof. (See Miescke (1979)).

With this lemma, it is easy to get

$$p_i(x) = P(\mu_i = u[k] | \underline{x})$$

$$= \int I_{\{\mu_i = u[k]\}} d\nu \left(\left(\frac{Y}{q+\gamma} \right) x, \frac{\gamma q}{q+\gamma} I \right)^{(\mu)}$$

where $\nu_{(\mu, V)}$ is the normal distribution with mean $\underline{\mu}$ and variance-covariance matrix V .

We can rewrite $p_i(x)$ as

$$p_i(x) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi(t + (\frac{\gamma}{q(q+\gamma)})^{\frac{1}{2}}(x_i - x_j)) d\nu(t).$$

Let $\gamma = \sigma_0^2$, $q = \sigma^2/n$, we have

$$p_i(x) = \int_{-\infty}^{\infty} \prod_{j \neq i} \phi(t + (\frac{\sigma_0^2}{2 \cdot \frac{\sigma^2}{n} + \sigma_0^2})^{\frac{1}{2}}(x_i - x_j)) d\nu(t).$$

The above expression is exactly the same as that of the independent prior
Case I, Ib.

Case V. Under normal assumption as before, but suppose σ_i 's are unknown
and that neither σ_i 's nor n_i 's are all equal.

Suppose we have no prior information about (μ, σ) , for each individual population τ_i assign prior $p(\mu_i, \sigma_i) \cdot \sigma_i^{-1}$ then we have (See Box and Tiao (1973)) that the posterior density of μ_i , given $x_i = x_i^+ (x_{i1}, \dots, x_{in_i})$ is

$$p(\mu_i | x_i) = \frac{(s_i / \sqrt{n_i})^{-1}}{B(\frac{v_i}{2}, \frac{1}{2}) \sqrt{v_i}} [1 + \frac{n_i (\mu_i - x_i)}{v_i s_i^2}]^{-\frac{v_i}{2} + 1}$$

where s_i^2 is the sample variance, $B(\cdot, \cdot)$ is a Beta function and $v_i = n_i - 1$.

Hence

$$p(t_i = \frac{\mu_i - x_i}{s_i / \sqrt{n_i}} | x_i) = \frac{1}{B(\frac{v_i}{2}, \frac{1}{2}) \sqrt{v_i}} (1 + \frac{t_i^2}{v_i})^{-\frac{v_i}{2} + 1}$$

which is the density of the student's t distribution with $v_i (n_i - 1)$ degrees of freedom.

Using this result we can write the formula of $p_i(x)$ by

$$\begin{aligned} p_i(x) &= P(\mu_i > u_j, \forall j \neq i | x) \\ &= \int_{j \neq i} T v_j \left(t \frac{s_i / \sqrt{n_i}}{s_j / \sqrt{n_j}} + \frac{x_i - x_j}{s_j / \sqrt{n_j}} \right) d T v_i(t) \end{aligned}$$

where $v_i = n_i - 1$, $i = 1, \dots, k$

$$s_i^2 = \sum_{r=1}^{n_i} (x_{ir} - x_i)^2$$

$T v_i$ is the c.d.f. of t distribution with v_i degrees of freedom.

When v_i 's are large, t distribution approaches normal distribution, hence, for large n_i , $i = 1, \dots, k$, we can replace T by Φ .

Case VI. Suppose we are interested in finding a subset which contains the population with the smallest variance; i.e., we define the best population as the one with the smallest variance, and suppose that we have no prior information about σ . In this case, it is reasonable to assume that

$$p(\mu, \sigma^2)^{-1}, \text{ if } \mu \text{ is unknown}$$

$$p(\mu)^{-1}, \text{ if } \mu \text{ is known.}$$

Let

$$v_j = n_j, v_j s_j^2 = \sum_{r=1}^k (x_{jr} - \bar{x}_j)^2 \text{ if } \mu \text{ is known}$$

$$v_j = n_j - 1, v_j s_j^2 = \sum_{r=1}^k (x_{jr} - \bar{x}_j)^2 \text{ if } \mu \text{ is unknown, } n_j \geq 1 \quad j = 1, 2, \dots, k$$

$$s^2 = (s_1^2, \dots, s_k^2), \bar{x} = (\bar{x}_{11}, \dots, \bar{x}_{1n_1}, \dots, \bar{x}_{kn_k})$$

and Y_{ij} be the random variable with c.d.f. $\chi_{v_j}^2$ which is the χ^2 distribution with v_j degrees of freedom.

Then for either case (μ known or unknown), we have

$$\begin{aligned} p_i(x) &= P(\sigma_i^2 = s_{[1]}^2 | \bar{x} = x) \\ &= P(\sigma_i^2 \leq v_j^2, \forall j \neq i | \bar{x} = x) \\ &= P\left(\frac{v_j s_j^2}{v_j^2} \leq \frac{\sigma_i^2}{v_i^2} \left(\frac{v_j s_j^2}{v_i^2}\right), \forall j \neq i | s^2 = s^2\right) \\ &= P\left(Y_{ij} \leq \frac{v_j s_j^2}{v_i^2} \left(\frac{v_j s_j^2}{v_i^2}\right), \forall j \neq i | s^2 = s^2\right) \\ &= \int_0^{\infty} \frac{v_j^2}{v_j - v_i^2} \left(\frac{v_j s_j^2}{v_i^2}\right) u^{v_j s_j^2 - 1} e^{-u/v_j} du \end{aligned}$$

$$= \int_0^{\infty} \sum_{j \neq i} \frac{s_j^2}{s_i^2} (u - \frac{s_j^2}{s_i^2}) d\psi(u) \quad \text{if } n_1 = \dots = n_k = 1.$$

With these $p_1(x), \dots, p_k(x)$ we can apply Bayes-P* rules ψ^B and ψ_{NR}^B easily.

Lemma 2.5.2. In Case VI, ψ^B and ψ_{NR}^B are just * a.e. and (scale) translation invariant.

* Here the definition of the "just" property for a selection rule is

$$\psi_i(s_i^2) < \psi_j(s_j^2) \text{ if } s_i^2 > s_j^2, s_j^2 < s_i^2, \forall j \neq i.$$

2.6. Comparison of Selection Rules ψ^B and ψ^M in the Normal Location

Parameter Case

We have k normal populations with a common known variance σ^2 and common sample size n . For this case Gupta (1956) proposed and studied the procedure ψ^M .

ψ^M : Select x_i iff $x_i > x_{[k]} + d \frac{\sigma}{\sqrt{n}}$, $i = 1, \dots, k$ where $d = d(\psi^M)$ is to be determined by

$$\inf_{\theta \in \Omega} P(CS|\psi^M) = p^*$$

and Ω is the parameter space.

We will show that $\psi^M \in \mathcal{L}_{NR}(\cdot, p^*)$ where \cdot is the locally uniform prior distribution. For fixed p^* and k , let d be determined by

$$\int_{-\infty}^{\infty} \psi^{k-1}(t + d) d\psi(t) = p^*. \quad (1.6.1)$$

Let

$\mathcal{X}_i = \{\text{all possible observed values}\} \subset \mathbb{R}^k$

$$\mathcal{X}_{i-1} = \{x \in \mathcal{X} | x_{[k]} - d \frac{\beta}{\sqrt{n}} < x_{[i]}\}$$

$$\mathcal{X}_i = \{x \in \mathcal{X} | x_{[i-1]} < x_{[k]} - d \frac{\beta}{\sqrt{n}} < x_{[i]}\}, \quad 2 \leq i \leq k$$

$$\mathcal{X}_i^{(1)} = \{x \in \mathcal{X} | x_{[1]} = x_{[i-1]} < x_{[k]} - d \frac{\beta}{\sqrt{n}} < x_{[i]}\} \subset \mathcal{X}_i$$

$$\mathcal{X}_i^{(2)} = \{x \in \mathcal{X} | x_{[1]} = x_{[i-1]} < x_{[k]} - d \frac{\beta}{\sqrt{n}} = x_{[i]} = x_{[k-1]}\} \subset \mathcal{X}_i^{(1)}$$

then we have the following theorem.

Theorem 2.6.1. Given a number $P^* (\frac{1}{k} < P^* < 1)$ and locally uniform prior for each population π_i , $x = \underline{x} \in \mathcal{X}_i$, then

$$P(\text{CS}_{i-1}^M, \underline{x} = \underline{x}) \geq q^*(i)$$

where

$$q^*(i) = \frac{k-i}{k-1} (1 - P^*) + P^*.$$

Hence

$$\underline{x} \in \mathcal{X}_{NR}(i, P^*) .$$

Proof. It is sufficient to show that

$$\inf_{x \in \mathcal{X}_i} \sum_{j=1}^k p_{ij} p_{ij}(x) = q^*(i) = \frac{k-i}{k-1} (1 - P^*) + P^*.$$

Since $\underline{x} \in \mathcal{X}_{i-1}$,

$$P(\text{CS}_{i-1}^M, \underline{x}) = \inf_{x \in \mathcal{X}_i} P(\text{CS}_{i-1}^M, x)$$

$$\inf_{x \in \mathcal{X}_{i-1}} \sum_{j=1}^k p_{ij} p_{ij}(x) .$$

Without loss of generality we can assume $\frac{c}{\sqrt{n}} = 1$.

Since

$$\sum_{j=1}^k p_{[j]}(x) = 1 \quad \forall x \in \mathcal{Z}, \text{ and } V = i$$

$p_{[i]}(x)$ is nonincreasing for all $x_{[j]}, j \leq i - 1$, we have

$$\begin{aligned} \inf_{x \in \mathcal{Z}_i} \sum_{j=i}^k p_{[j]}(x) &= \inf_{x \in \mathcal{Z}_i} (1) \sum_{j=i}^k p_{[j]}(x) \\ &= 1 - \sup_{x \in \mathcal{Z}_{\bar{j}}} (1) \sum_{j=1}^{i-1} p_{[j]}(x) \\ &= 1 - \sup_{x \in \mathcal{Z}_{\bar{j}}} (1) \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \sum_{\substack{t \geq 0 \\ j \neq i}} \psi(t + x_{[j]} - x_{[j]}) d\psi(t) \\ &= 1 - \sup_{x \in \mathcal{Z}_{\bar{j}}} (1) \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \sum_{\substack{t \geq 0 \\ j \neq i}} \psi(t + x_{[j]} - x_{[j]}) \\ &\quad \cdot \sum_{\substack{t \geq 0 \\ j \neq i}} \psi(t + x_{[j]} - x_{[j]}) d\psi(t) \\ &= 1 - \sup_{x \in \mathcal{Z}_{\bar{j}}} (1) \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \sum_{\substack{t \geq 0 \\ j \neq i}} \psi(t + x_{[j]} - x_{[j]}) \\ &\quad \cdot \psi^{i-2}(t) d\psi(t) \\ &= 1 - \sum_{j=1}^{i-1} \int_{-\infty}^{\infty} \psi(t - d) \psi^{i-2}(t) d\psi(t) \quad (2.6.7) \end{aligned}$$

$$\begin{aligned}
&= 1 - (i-1) \int_{-\infty}^{\omega} \phi(t-d)\psi^{k-2}(t)d\phi(t) \\
&= (k-i) \int_{-\infty}^{\omega} \phi^{k-2}(t)\phi(t-d)d\phi(t) \\
&\quad + \int_{-\infty}^{\omega} \phi^{k-1}(t+d)d\phi(t) \tag{2.6.3}
\end{aligned}$$

The superimum of (2.6.3) occurs when $x \in \mathcal{X}_i^{(2)}$. The last equality follows from the identity

$$\begin{aligned}
&(k-i) \int \phi^{k-2}(t)\phi(t-d)d\phi(t) \\
&= 1 - \int \phi^{k-1}(t+d)d\phi(t),
\end{aligned}$$

which can be shown by the integration by parts. By (2.6.1), the second term of (2.6.3) equals P^* ; then use the integration by parts to the first term of (2.6.2), we get

$$\begin{aligned}
\inf_{x \in \mathcal{X}_i} \sum_{j=i}^k p_{[j]}(x) &= \frac{k-i}{k-1} [1 - P^*] + P^* \tag{2.6.4} \\
&= q^*(i).
\end{aligned}$$

Remark 2.6.1. If the procedure ψ^M selects $x_{(k)}$ only, i.e. $X = x \in \mathcal{X}_k$, then by Theorem 2.6.1 we have $p_{[k]}(x) \geq P^*$ so that ψ^B or ψ_{NR}^B selects $x_{(k)}$ only. But the converse is not necessarily true.

Remark 2.6.2. For the case $k = 2$, $\psi_{NR}^B = \psi^M$ a.e. For any given $X = x$: if $x \in \mathcal{X}_2$, then $p_{[2]}(x) \geq P^*$, hence ψ^M and ψ_{NR}^B select the population $x_{(2)}$ associated $x_{[2]}$. If $x \in \mathcal{X}_1$, and $x_{[2]} - d \frac{\sigma}{\sqrt{n}} < x_{[1]}$ then ψ^M and

ψ_{NR}^B select both populations π_1 and π_2 . Since

$$P(X_{[2]} - d \frac{\sigma}{\sqrt{n}} = X_{[1]}) = 0,$$

we have $\psi_{NR}^B = \psi^M$ a.e. .

Remark 2.6.3. The above Theorem and Remark 2.6.1 gives us a lower

bound on the value of $\sum_{i=1}^k P_{[i]}(x)$, over all $x \in \cup_i \pi_i$. The exact value of the difference of the selected sizes between ψ^M and ψ^B depends on the observations.

2.7. Applications to Select $\max_{1 \leq i \leq k} \pi_i$, $\rho_i = \frac{\mu_i - a}{\sigma_i}$ for Normal

Distribution $N(\mu_i, \sigma_i^2)$, $i = 1, \dots, k$

Let π_1, \dots, π_k be k independent normal populations with mean μ_i and variance σ_i^2 , both μ_i and σ_i are unknown. For the goal of finding a

random subset which contains the population with maximum $\rho_i = \frac{\mu_i - a}{\sigma_i}$

for some given constant a , we assume that apriori (μ_i, σ_i) , $i = 1, \dots, k$ are independent. Suppose we have n_i independent observations

x_{i1}, \dots, x_{in_i} from π_i , and let X_i be their sample mean, $i = 1, \dots, k$.

Let y_1, \dots, y_n be i.i.d. $\sim N(\mu, \sigma^2)$. If no prior information is available to (μ, σ) , we could assign a locally uniform prior $p(\mu, \sigma) \propto \sigma^{-1}$ to (μ, σ) , (see Box and Tiao (1973)). And the posterior joint distribution of $\mu' = \mu - a$ and σ , given observations

$Y = y = (y_1, \dots, y_n)$ is given by

$$P(\mu^*, \sigma^* | y) = k \sigma^{-(n+1)} \exp \left\{ -\frac{1}{2\sigma^2} [vs^2 + n(y^* - \bar{y})^2] \right\}$$

where

$$\left. \begin{aligned} y^* &= y - a, \quad \bar{y} = \sum_{i=1}^n y_i/n \\ vs^2 &= \sum (y_i - \bar{y})^2, \quad v = n - 1 \\ k &= \sqrt{\frac{n}{2\pi}} \left[\frac{1}{2} \Gamma\left(\frac{v}{2}\right) \right]^{-1} \left(\frac{vs}{2} \right)^{\frac{v}{2}}. \end{aligned} \right\} \quad (2.7.1)$$

Let $t = \sqrt{n}(\mu - a)/\sigma$, with (2.7.1) the posterior distribution of t , given $Y = y$ is

$$p(t | Y = y) = p(t | t)$$

$$= \{2^{\frac{v}{2}-1} \Gamma\left(\frac{v}{2}\right)\}^{-1} \left(\frac{v}{v+t^2}\right)^{\frac{v}{2}} \exp \left\{ -\frac{1}{2} \left(\frac{vt}{v+t^2} \right)^2 \right\} \Gamma(v) I_{v-1} \left(\frac{vt}{v+t^2} \right)^{1/2}$$

where

$$t = \sqrt{n}(y - a)/s, \quad v = n - 1$$

$$I_v(x) = \int_0^\infty (2\pi \Gamma(v))^{-1} u^v \exp \left\{ -\frac{1}{2} (u + x)^2 \right\} du.$$

Now, let $p(\mu_i, \sigma_i) = \sigma_i^{-1}$ be the assigned locally uniform prior to (μ_i, σ_i) . Then let $x = (x_{i1}, \dots, x_{1n_1}, \dots, x_{kn_k})$, we have

$$p_i(x) = P(\mu_i \in [\cdot, \cdot] | x)$$

$$= P\left(\frac{\mu_i - a}{\sigma_i} \leq \max_{1 \leq j \leq k} \left(\frac{\mu_j - a}{\sigma_j} \right) | x\right)$$

$$\begin{aligned}
 &= P\left(\sqrt{\frac{n_j}{n_i}} \xi_j > \xi_i \mid t\right) \\
 &= \int_{j \neq i} G_{\xi_j}\left(\sqrt{\frac{n_j}{n_i}} z \mid t\right) d G_{\xi_i}(z \mid t) \quad (2.7.1) \\
 &= \int_{j \neq i} G_{\xi_j}(z \mid t) d G_{\xi_i}(z \mid t) \quad \text{if } n_1 = \dots = n_k = n,
 \end{aligned}$$

where G_{ξ_j} is the posterior c.d.f. of ξ_j , given x or t .

By (2.7.2), the Bayes-P* procedure is completely specified.

If the prior distribution for (μ, σ) is the conjugate distribution (see Raiffa and Schlaifer (1960)), then

$$\begin{aligned}
 p(\mu, \sigma) &\propto \exp\left(-\frac{1}{2\sigma^2} n'(\mu - m')^2 + \frac{1}{\sigma} \cdot \exp\left(-\frac{1}{2\sigma^2} (\nu' v')\right)\right)^{-1/2} \\
 &= p(\mu|\sigma) p(\sigma)
 \end{aligned}$$

that is

$$p(\mu|\sigma) \sim N(m', \sigma^2/n'), n' > 0$$

$$p(\sigma) \sim \frac{\nu' v'}{2} \cdot \sigma^{-2}, \nu', v' > 0.$$

Let

$$x' = \frac{nx + n'm'}{n + n'}, x \text{ is the sample mean.}$$

$$u^2 = [(n-1)s^2 + v'v' + ((nn')/(n+n'))(x - m')^2]/ *$$

$$v^* = (n-1) + v' + 1$$

$$\tau^* = (n + n')^{1/2}(\mu - a)/u$$

$$t^* = (n + n')^{1/2}(x' - a)/u,$$

the posterior distribution of τ^* , given x is $p(\tau^*|x) \sim p(\tau^*|t^*)$ which has the same form as $p(\tau|t)$, but replace τ , t , a by τ^* , t^* , a^* .

thus for the conjugate prior case, we get

$$\begin{aligned} p_i(\cdot) &= p(\cdot | \beta_i^*, \{\epsilon\})^{-1} \\ &\propto \int_{\beta \neq \beta_i^*} G_{\beta_i^*}(z|t^*) dG_{\beta_i^*}(z|t^*) \quad (\text{from (2.7.2)}) \\ &\propto \int_{\beta \neq \beta_i^*} G_{\beta_i^*}(z|t^*) dG_{\beta_i^*}(z|t^*) \quad (\text{if } n_i = 1) \end{aligned}$$

where $G_{\beta_i^*}$ is the posterior c.d.f. of β_i^* given x or t .

Note that (2.7.3) has the same form as (2.7.2), but replace β, t by β^*, t^* .

2.8. Applications to Poisson Distributions and Poisson Process

2.8.1. Poisson Distributions Case

Suppose that $\gamma_1, \dots, \gamma_k$ are k independent Poisson populations, where the independent observations x_{i1}, \dots, x_{in_i} from γ_i have the Poisson density with parameter γ_i ; denoted by $P(\cdot | \gamma_i)$, $i = 1, \dots, k$.

Let y_1, \dots, y_n be i.i.d. with $p(\cdot | \gamma)$. If we use non-informative prior $p(\cdot) \propto \gamma^{-1/2}$ (Box and Tiao (1973)), then given $y = y = (y_1, \dots, y_n)$, we have the posterior density as follows:

$$p(\cdot | y) = e^{ny - \frac{1}{2}} / \exp(-n\gamma)$$

where

$$y = \frac{1}{n} \sum_{i=1}^n y_i \quad \text{and} \quad \gamma = n(ny + \frac{1}{2}) / [ny + \frac{1}{2}]^{-1}$$

We see that $2n_i \chi_i^2 \sim \chi_{2n_i+1}^2$, the chi-square distribution with $2n_i+1$ degrees of freedom. Hence by using non-informative prior

$p(\chi_i) = \chi_i^{-1/2}$ for each population χ_i , we have

$$\begin{aligned} p_i(x) &= p(\chi_i = \chi_{[k]} | x) \\ &= \int_0^\infty \prod_{j \neq i} \chi_j^2(z) \frac{n_j}{\chi_j^2(z)} d\chi_j^2(z) \end{aligned}$$

where

$$\chi_i = 2n_i x_i + 1, \quad x_i = \sum_{j=1}^{n_i} x_{ij}/n_i.$$

If $n_1 = \dots = n_k$, then

$$p_i(x) = \int_0^\infty \prod_{j \neq i} \chi_j^2(z) \frac{n_j}{\chi_j^2(z)} d\chi_j^2(z).$$

With $p_i(x)$, $i = 1, \dots, k$, we can apply Bayes-P* selection rules

χ_i^B and χ_{NR}^B easily to select a subset which contains the population with the largest parameter χ . On the other hand, if we are interested in selecting the population with the smallest parameter χ , then

$$\begin{aligned} p_i(x) &= \int_0^\infty [1 - \chi_i^2(z)] \frac{n_i}{\chi_i^2(z)} d\chi_i^2(z) \\ &= \int_0^\infty [1 - \chi_i^2(z)] \frac{n_i}{\chi_i^2(z)} d\chi_i^2(z) \quad \text{if } n_1 = \dots = n_k. \end{aligned}$$

In this case, the simulation results for selection procedures and χ_{NR}^B are tabulated on Table VII.

2.8.2. Poisson Processes Case

Suppose we have k independent Poisson processes

$\{X^{(1)}(t)\}, \dots, \{X^{(k)}(t)\}$ with expected arrival times equal to

$\frac{1}{\lambda_1}, \dots, \frac{1}{\lambda_k}$, respectively. Hence for the processes $\{X^{(i)}(t)\}$, the

probability that there are m_i arrivals until time t_i is

$$p(X^{(i)}(t_i) = m_i | \lambda_i, t_i) = \frac{(t_i \lambda_i)^{m_i}}{m_i!} e^{-t_i \lambda_i}$$

$$i = 1, 2, \dots, k, \quad m_i = 0, 1, 2, \dots$$

If there exists no prior information, then we use the non-informative prior $p(\lambda_i) \propto \lambda_i^{-1/2}$ for all processes. Therefore, we get the posterior density function of λ_i , given (m_i, t_i) as follows:

$$\begin{aligned} p(\lambda_i | X^{(i)}(t_i) = m_i, t_i) &= p(\lambda_i | m_i, t_i) \\ &= \frac{(t_i \lambda_i)^{m_i + \frac{1}{2} - 1}}{\Gamma(m_i + \frac{1}{2})} t_i e^{-t_i \lambda_i}. \end{aligned}$$

Thus $2t_i \lambda_i$ has χ^2 distribution with $2m_i + 1$ degrees of freedom, given the number m_i of arrivals before time t_i .

Let $m = (m_1, \dots, m_k)$ and $t = (t_1, \dots, t_k)$, then it can be shown that the Poisson process $\{X^{(i)}(t)\}$ has the maximum parameter (or minimum mean waiting time), given (m, t) is

$$p_i(m, t) = \int_0^\infty \prod_{j \neq i} x_{2m_j+1}^2(y) \frac{t_j}{t_i} d x_{2m_i+1}^2(y) \quad i = 1, \dots, k. \quad (7.2.1)$$

Here we list two special cases which are of interest.

- (a) Observations of all processes are obtained in a common time interval $[s_i, t + s_i]$. Since Poisson process is stationary, we can assume that $s_i = 0$, and $t_1 = \dots = t_k = t$. In this case

$$p_i(m, t) = \int_0^\infty \prod_{j \neq i} x_{2m_j+1}^2(y) d x_{2m_i+1}^2(y)$$

which is independent of t .

- (b) All m_i 's are equal, i.e. we fix m first, then net observations t . Hence

$$p_i(m, t) = \int_0^\infty \prod_{j \neq i} x_{2m+1}^2(y) \frac{t_j}{t_i} d x_{2m+1}^2(y).$$

There is an alternative way to approach the cases (a) and (b). Let T_i be the waiting time of the n th arrival in the i th process, then T_i has a gamma distribution with density given by

$$p(t) = \frac{\lambda_i^{m_i}}{\Gamma(m_i)} (\lambda_i t)^{m_i-1} e^{-\lambda_i t} \quad t > 0.$$

If we have only non-informative prior $p(\lambda) \propto \lambda^{-1/2}$; then, given m_i and t_i , $\lambda_i t_i$ has posterior distribution $\frac{?}{x_{2m_i+1}^2}$, therefore the formula of $p_i(m, t)$ we get here is exactly the same as before.

Remark 2.8.1. Under non-informative prior, in comparing the subset selection problem in k Poisson distributions with the problem in k Poisson processes, it is easily seen that Poisson distributions model is a special case of Poisson processes model, namely, $t_i = n_i$ where n_i denotes the sample size of the i th Poisson population.

2.8.3. Relation Between Selection from Poisson Processes and Selection from Populations with Gamma or Exponential Distribution

Suppose we have k independent populations, the i th population having the gamma distribution with parameters $\alpha = m_i$ (known), $\beta = 1/\lambda_i$ (unknown). Since the random variable T_i , the waiting time until m_i arrivals in a Poisson process with parameter λ_i , has a gamma distribution with parameters $\alpha = m_i$, $\beta = 1/\lambda_i$. If the m_i 's are given and if the goals for both selection problems are the same, namely, to select a subset containing the population (process) with the largest parameter λ , then it is easily seen that these are identical problems. Note that in the selection problem of Poisson processes, m_i 's might not be the preassigned values but are given random observations whenever t_i 's are preassigned values. In this case, the selection problem of Poisson processes is different from that of the gamma distributions.

If the process associated with the minimum parameter λ (or the maximum waiting time) is the best, then the posterior probability of process $x^{(i)}(t)$ to be the best is analogous to the one obtained

before with the modifications that the integrand function

$$\sum_{j \neq i}^n x_{2m_j+1}^2 (y \frac{t_j}{t_i})$$

of (2.8.1) is replaced by

$$\sum_{j \neq i}^n [1 - x_{2m_j+1}^2 (y \frac{t_j}{t_i})].$$

2.9. Comparison of the Performance of ψ^B , ψ_{NR}^B , ψ^M and ψ^{MED}

Let π_i , $i = 1, \dots, k$ be k independent populations, where π_i has the associated c.d.f. $F(x, \theta_i) = F(x - \theta_i)$ with unknown location parameter θ_i . Let $f(x, \theta_i) = f(x - \theta_i)$ be the p.d.f. The goal is to find a small (nontrivial) subset which contains the best.

The following subset selection procedure, ψ^{MED} based on sample medians is due to Gupta and Singh (1980).

ψ^{MED} : Select π_i if and only if $X_i \leq \bar{X}_{[k]} - d$

where \bar{X}_i is the median of the $2m+1$ random observations from population π_i and $\bar{X}_{[k]} = \max_i \bar{X}_i$. The value d is determined by the following equation so that the P^* -condition is met.

$$\int_{-\infty}^{\infty} G(u + d)^{k-1} g(u) du = P^*$$

where

$$g(u) = \frac{(2m+1)!}{(m!)^2} [F(u)]^m [1 - F(u)]^m f(u)$$

$$G(u) = I_F(u)^{(m+1, m+1)}$$

$I_y(p, q)$ is the incomplete beta function.

In this section we use Monte Carlo simulation techniques to compare the performance of selection procedures ψ^B , ψ_{NR}^B , ψ^M and ψ^{MED} in the normal means problem. Because both rules ψ^M and ψ^{MED} are not based on any prior information about the unknown parameters, we assume that the prior distribution π for both ψ^B and ψ_{NR}^B is locally uniformly distributed. Since the selection procedure ψ^M satisfies both the P*-condition and the posterior-P* condition wrt the locally uniform priors, it makes sense to compare the Bayes-P* procedures ψ^B and ψ_{NR}^B with ψ^M and compare ψ^M with ψ^{MED} in terms of efficiency which is the ratio of the probability of a correct selection to the expected selected size. For studying the robustness of these four rules, ψ^B , ψ_{NR}^B , ψ^M and ψ^{MED} , we change the true distribution to non-normal distributions, namely, the logistic, Laplace (the double exponential) and the gross error model (the contaminated distribution), but keep the selection procedure unchanged (i.e. still based on the normal assumption). The Monte Carlo simulation results for both equal distances of the parameters and slippage cases are tabulated. In the simulation study all generated random variables are adjusted to have variance 1. Each time we generate five random variables with the given distribution of each population, then apply the selection procedures. The simulation process is repeated 100 times for each random variable. The relative frequency of selecting the population π_1 is

used as an approximation to the probability of selecting the population π_i . The sum of relative frequency of selecting each population π_i , $i = 1, \dots, k$ is treated as an approximation of the expected selected size. The efficiency EFF of each selection procedure is approximated by the ratio of relative frequency of selecting the best one to the expected size. The simulation results indicate that in all cases we have the performance

$$\phi_B > \phi_{NP} > \phi_M.$$

It should be noted that in the above comparison of the performances, we restrict attention to these rules which satisfy the posterior P* condition. For small sample size, the efficiency of rule ϕ_M tends to be larger than ϕ_{MED} under P*-condition.

Remark 2.9.1. The Laplace distribution has the density function

$$f(x - \mu) = \frac{1}{2} e^{-|x-\mu|}, \quad x \in \mathbb{R}$$

for which the variance is 2.

The logistic distribution has the density function

$$f(x - \mu) = \frac{e^{-(x-\mu)}}{1 + e^{-(x-\mu)}}^2$$

for which the variance $\text{Var}(X) = \frac{\pi^2}{3}$.

The gross error model we used has the density function

$$f(x - \mu) = (1 - \beta) f(x - \mu + \beta) \cdot \left(\frac{\lambda}{\beta}\right)^{\lambda}, \quad \lambda > 0, \beta > 0$$

for which $f(\cdot)$ is the p.d.f. of $N(0,1)$ and the variance $\text{Var}(X) = (1 - \beta) + \beta^2 + \beta^2 = 3.25$.

The efficiency of a selection procedure π is defined by

$$EFF(\pi) = \frac{P_{\pi}(S^*)}{E_{\pi}(S^*)}$$

where $E_{\pi}(S^*)$ is the expected selected size.

Discussion and Conclusion

For Table VIII.1 and Table VIII.2 (equal distances case) the P^* is .99 and .90 respectively, the common sample size $n = 5$, $k = 5$. If the k populations have normal distributions with the unknown parameter configuration $(\mu_1, \dots, \mu_k; \sigma^2)$, common variance 1. From both tables the performance based on either the efficiency or the expected selected size is

$$\frac{P}{P^*} > \frac{S}{NS} > \frac{M}{MD}$$

If the posterior- P^* condition is considered, and

$$M > MD$$

under the P^* -condition,

when the true distributions are not normal, but the logistic, the Laplace or the trias error model, the results are very close to the normal case, none of the four rules are robust. From Table VIII.3 different P^* 's are three times than the corresponding ones in Table VIII.1. This is to be expected because the value of P^* is smaller in the second table.

In Table VIII.1 and VIII.2 therefore under the P^* rule, 99 and 90% respectively, the common sample sizes of k are 5 for the k populations.

have normal distributions with unknown parameter configurations (μ_1, \dots, μ_k) , common variance 1. From both tables the performance based on either the efficiency or the expected selected size is:

$$\psi^B \geq \psi_{NR}^B \geq \psi^M$$

if the posterior-P* condition is considered, and if $\sqrt{n} = 1$

$$\psi^M > \psi^{MED}$$

under the P*-condition.

Note that in both equal distances and slippage cases when $\psi^M = 1$, that is the population means are not very close, the procedures ψ^B and ψ_{NR}^B , wrt the locally uniform priors, always satisfy not only the posterior-P* condition but also $P(\text{CS}|\psi^B \text{ or } \psi_{NR}^B) \geq P^*$, and the expected selected size of the selection procedure ψ^B or ψ_{NR}^B is much less than the selection procedures ψ^M and ψ^{MED} . For example, in the normal equal distances case, $P^* = .99$, $k = 5$, $\sqrt{n} = 4$,

$$E(S|\psi^{MED}) - E(S|\psi_{NR}^B) = 0.382;$$

in the normal slippage case, $P^* = .99$, $k = 5$, $\sqrt{n} = 4$

$$E(S|\psi^{MED}) - E(S|\psi_{NR}^B) = 1.560.$$

TABLE III

For procedures \hat{B} and \hat{E} and the parameter configurations $(.5, .5, .5, .5)$ of k Poisson populations, this table gives the values (based on simulation) of the probability of selecting the population with parameter $.5i$, $i=1, \dots, k$, and the expected selected size ES . The prior distribution for each population is $B(\lambda_i, \lambda_i)$.

 $n = 10$

k	P*				0.75			
	0.99	0.95	0.90	0.75	0.99	0.95	0.90	0.75
1	.993	1.000	.966	.950	.990	.950	.946	.990
2	.670	.820	.450	.670	.324	.490	.110	.190
3	1.663	1.820	1.415	1.660	1.274	1.480	1.055	1.180
4	.968	1.000	.990	1.000	.967	1.000	.929	1.000
5	.741	.850	.388	.630	.321	.510	.171	.280
6	2.05	.330	.120	.130	.065	.110	.021	.050
7	1.944	2.180	1.498	1.760	1.354	1.620	1.121	1.336
8	.993	1.060	.981	1.000	.966	.990	.911	.950
9	.745	.829	.560	.730	.259	.400	.175	.290
10	3.69	.496	.070	.140	.067	.100	.025	.050
11	.675	.110	.041	.060	.024	.046	0	0
12	2.167	2.426	1.653	1.430	1.316	1.530	1.112	1.290
13	.996	1.090	.981	.990	.985	1.000	.950	.980
14	.737	.850	.503	.650	.367	.550	.128	.260
15	3.55	.479	.102	.160	.133	.210	.013	.050
16	.667	.090	.015	.030	.015	.020	.006	.010
17	2.154	.413	.006	.016	0	0	0	0
18	.997	1.154	1.667	1.650	1.498	1.760	1.347	1.460

TABLE VIII. 1

Efficiency (EFF) and Expected Selected Size (ES) (based on simulation) of B_M , B_{NR} , and B_{MED} when the unknown means of the k populations are $\beta_1, \dots, \beta_k + (k-1)\delta$; the common variance is 1, common sample size $n = 5$ and the prior for β and β_{NR} is locally uniformly distributed.

		Normal		Logistic		Tage		Cross error	
		Eff	ES	Eff	ES	Eff	ES	Eff	ES
$n = 3$	B	.254	3.809	.250	3.963	.259	3.772	.248	3.956
	NR	.238	4.110	.233	4.290	.238	4.120	.230	4.301
	M	.208	4.810	.207	4.840	.210	4.720	.207	4.840
	MED	.208	4.810	.202	4.940	.202	4.940	.201	4.930
$n = 5$	B	.333	2.977	.332	3.005	.336	2.941	.329	3.032
	NR	.305	3.280	.304	3.290	.302	3.280	.304	3.291
	M	.250	4.000	.234	4.280	.246	4.030	.233	4.240
	MED	.232	4.310	.224	4.460	.217	4.600	.214	4.670
$n = 7$	B	.541	1.847	.541	1.829	.541	1.884	.559	1.779
	NR	.486	1.956	.481	2.085	.498	2.010	.510	1.960
	M	.417	2.400	.417	2.400	.515	2.410	.437	2.290
	MED	.369	2.720	.351	2.350	.367	2.76	.369	2.720
$n = 9$	B	.625	1.212	.955	1.162	.621	1.217	.865	1.156
	NR	.730	1.370	.666	1.240	.746	1.340	.695	1.250
	M	.676	1.420	.694	1.420	.630	1.470	.671	1.470
	MED	.676	1.720	.576	1.730	.671	1.730	.671	1.730

TABLE VIII. 2

Efficiency (EFF) and Expected Selected Size (ES) (based on simulation) of B , B_{NR} , M and MED when the unknown means of the k populations are $\dots, \dots, \dots + (k-1)\cdot$; the common variance is 1, common sample size $n = 5$ and the prior for B and B_{NR} is locally uniformly distributed.

TABLE VI

μ	σ^2	$\hat{\mu}$	$\hat{\sigma}^2$	bias	MSE	Gross error
0.000	0.000	0.000	0.000	-0.000	0.000	0.000
0.000	0.001	0.000	0.001	-0.001	0.001	0.001
0.000	0.002	0.000	0.002	-0.002	0.002	0.002
0.000	0.003	0.000	0.003	-0.003	0.003	0.003
0.000	0.004	0.000	0.004	-0.004	0.004	0.004
0.000	0.005	0.000	0.005	-0.005	0.005	0.005
0.000	0.006	0.000	0.006	-0.006	0.006	0.006
0.000	0.007	0.000	0.007	-0.007	0.007	0.007
0.000	0.008	0.000	0.008	-0.008	0.008	0.008
0.000	0.009	0.000	0.009	-0.009	0.009	0.009
0.000	0.010	0.000	0.010	-0.010	0.010	0.010
0.000	0.011	0.000	0.011	-0.011	0.011	0.011
0.000	0.012	0.000	0.012	-0.012	0.012	0.012
0.000	0.013	0.000	0.013	-0.013	0.013	0.013
0.000	0.014	0.000	0.014	-0.014	0.014	0.014
0.000	0.015	0.000	0.015	-0.015	0.015	0.015
0.000	0.016	0.000	0.016	-0.016	0.016	0.016
0.000	0.017	0.000	0.017	-0.017	0.017	0.017
0.000	0.018	0.000	0.018	-0.018	0.018	0.018
0.000	0.019	0.000	0.019	-0.019	0.019	0.019
0.000	0.020	0.000	0.020	-0.020	0.020	0.020
0.000	0.021	0.000	0.021	-0.021	0.021	0.021
0.000	0.022	0.000	0.022	-0.022	0.022	0.022
0.000	0.023	0.000	0.023	-0.023	0.023	0.023
0.000	0.024	0.000	0.024	-0.024	0.024	0.024
0.000	0.025	0.000	0.025	-0.025	0.025	0.025
0.000	0.026	0.000	0.026	-0.026	0.026	0.026
0.000	0.027	0.000	0.027	-0.027	0.027	0.027
0.000	0.028	0.000	0.028	-0.028	0.028	0.028
0.000	0.029	0.000	0.029	-0.029	0.029	0.029
0.000	0.030	0.000	0.030	-0.030	0.030	0.030
0.000	0.031	0.000	0.031	-0.031	0.031	0.031
0.000	0.032	0.000	0.032	-0.032	0.032	0.032
0.000	0.033	0.000	0.033	-0.033	0.033	0.033
0.000	0.034	0.000	0.034	-0.034	0.034	0.034
0.000	0.035	0.000	0.035	-0.035	0.035	0.035
0.000	0.036	0.000	0.036	-0.036	0.036	0.036
0.000	0.037	0.000	0.037	-0.037	0.037	0.037
0.000	0.038	0.000	0.038	-0.038	0.038	0.038
0.000	0.039	0.000	0.039	-0.039	0.039	0.039
0.000	0.040	0.000	0.040	-0.040	0.040	0.040
0.000	0.041	0.000	0.041	-0.041	0.041	0.041
0.000	0.042	0.000	0.042	-0.042	0.042	0.042
0.000	0.043	0.000	0.043	-0.043	0.043	0.043
0.000	0.044	0.000	0.044	-0.044	0.044	0.044
0.000	0.045	0.000	0.045	-0.045	0.045	0.045
0.000	0.046	0.000	0.046	-0.046	0.046	0.046
0.000	0.047	0.000	0.047	-0.047	0.047	0.047
0.000	0.048	0.000	0.048	-0.048	0.048	0.048
0.000	0.049	0.000	0.049	-0.049	0.049	0.049
0.000	0.050	0.000	0.050	-0.050	0.050	0.050
0.000	0.051	0.000	0.051	-0.051	0.051	0.051
0.000	0.052	0.000	0.052	-0.052	0.052	0.052
0.000	0.053	0.000	0.053	-0.053	0.053	0.053
0.000	0.054	0.000	0.054	-0.054	0.054	0.054
0.000	0.055	0.000	0.055	-0.055	0.055	0.055
0.000	0.056	0.000	0.056	-0.056	0.056	0.056
0.000	0.057	0.000	0.057	-0.057	0.057	0.057
0.000	0.058	0.000	0.058	-0.058	0.058	0.058
0.000	0.059	0.000	0.059	-0.059	0.059	0.059
0.000	0.060	0.000	0.060	-0.060	0.060	0.060
0.000	0.061	0.000	0.061	-0.061	0.061	0.061
0.000	0.062	0.000	0.062	-0.062	0.062	0.062
0.000	0.063	0.000	0.063	-0.063	0.063	0.063
0.000	0.064	0.000	0.064	-0.064	0.064	0.064
0.000	0.065	0.000	0.065	-0.065	0.065	0.065
0.000	0.066	0.000	0.066	-0.066	0.066	0.066
0.000	0.067	0.000	0.067	-0.067	0.067	0.067
0.000	0.068	0.000	0.068	-0.068	0.068	0.068
0.000	0.069	0.000	0.069	-0.069	0.069	0.069
0.000	0.070	0.000	0.070	-0.070	0.070	0.070
0.000	0.071	0.000	0.071	-0.071	0.071	0.071
0.000	0.072	0.000	0.072	-0.072	0.072	0.072
0.000	0.073	0.000	0.073	-0.073	0.073	0.073
0.000	0.074	0.000	0.074	-0.074	0.074	0.074
0.000	0.075	0.000	0.075	-0.075	0.075	0.075
0.000	0.076	0.000	0.076	-0.076	0.076	0.076
0.000	0.077	0.000	0.077	-0.077	0.077	0.077
0.000	0.078	0.000	0.078	-0.078	0.078	0.078
0.000	0.079	0.000	0.079	-0.079	0.079	0.079
0.000	0.080	0.000	0.080	-0.080	0.080	0.080
0.000	0.081	0.000	0.081	-0.081	0.081	0.081
0.000	0.082	0.000	0.082	-0.082	0.082	0.082
0.000	0.083	0.000	0.083	-0.083	0.083	0.083
0.000	0.084	0.000	0.084	-0.084	0.084	0.084
0.000	0.085	0.000	0.085	-0.085	0.085	0.085
0.000	0.086	0.000	0.086	-0.086	0.086	0.086
0.000	0.087	0.000	0.087	-0.087	0.087	0.087
0.000	0.088	0.000	0.088	-0.088	0.088	0.088
0.000	0.089	0.000	0.089	-0.089	0.089	0.089
0.000	0.090	0.000	0.090	-0.090	0.090	0.090
0.000	0.091	0.000	0.091	-0.091	0.091	0.091
0.000	0.092	0.000	0.092	-0.092	0.092	0.092
0.000	0.093	0.000	0.093	-0.093	0.093	0.093
0.000	0.094	0.000	0.094	-0.094	0.094	0.094
0.000	0.095	0.000	0.095	-0.095	0.095	0.095
0.000	0.096	0.000	0.096	-0.096	0.096	0.096
0.000	0.097	0.000	0.097	-0.097	0.097	0.097
0.000	0.098	0.000	0.098	-0.098	0.098	0.098
0.000	0.099	0.000	0.099	-0.099	0.099	0.099
0.000	0.100	0.000	0.100	-0.100	0.100	0.100

TABLE IX. 1

Estimated average bias and standard deviation (ES) (based on simulation) of B_{M} , B_{NR} , \bar{B} and MED when the true variance of error term in simulations are 1, 1.5, 2, 3, 4; the common variance is 1, common sample size $n = 5$ and type prior for B and B_{P} is locally uniformly distributed.

TABLE IV. 1
 $\alpha = .5$,
 $P^* = .25$

n	β	Normal		Logistic		Laplace		Gauss	
		EFF	ES	EFF	ES	EFF	ES	EFF	ES
2	.2	4.132	.218	4.253	.226	4.203	.220	4.361	
3	.2	4.527	.212	4.580	.214	4.530	.212	4.670	
4	.2	4.373	.202	4.950	.200	4.900	.202	4.890	
5	.2	4.920	.202	4.940	.202	4.960	.200	5.000	
6	.2	4.137	.240	4.153	.240	4.062	.231	4.264	
7	.2	4.460	.224	4.460	.223	4.430	.216	4.580	
8	.2	4.461	.203	4.930	.203	4.670	.214	4.900	
9	.2	4.373	.201	4.970	.200	4.980	.200	5.000	
10	.2	4.526	.274	3.643	.263	3.701	.265	3.381	
11	.2	4.972	.253	3.360	.242	4.090	.269	3.720	
12	.2	4.626	.211	4.750	.214	4.680	.212	4.510	
13	.2	4.250	.207	4.510	.203	4.570	.207	4.270	
14	.2	4.746	.251	3.520	.251	3.520	.251	3.550	
15	.2	4.250	.222	4.250	.222	4.250	.222	4.250	
16	.2	4.746	.251	3.520	.251	3.520	.251	3.550	
17	.2	4.250	.222	4.250	.222	4.250	.222	4.250	
18	.2	4.746	.251	3.520	.251	3.520	.251	3.550	
19	.2	4.250	.222	4.250	.222	4.250	.222	4.250	
20	.2	4.746	.251	3.520	.251	3.520	.251	3.550	

TABLE IX. 2

Estimated average and standard selected size (SES) (based on simulation) of B_{B} , B_{M} , M_{B} and M_{M} when error variances in the populations are 1, 10, 100, 1000; the common variance is 1, common sample size is 100 and there are 1000, 500, 250 and 100 initially uninformative variables.

N	M	Logistic	Gaussian	Estimate		Gross error	
				EF	ES	EF	ES
10	10	1.250	1.250	.238	2.734	.256	2.856
10	20	1.262	1.262	.228	3.160	.240	3.250
10	30	1.264	1.264	.211	4.400	.212	4.490
10	40	1.264	1.264	.202	4.840	.204	4.900
10	50	1.264	1.264	.194	5.354	.203	5.471
10	60	1.264	1.264	.186	5.905	.205	6.015
10	70	1.264	1.264	.178	6.500	.206	6.616
10	80	1.264	1.264	.170	7.150	.207	7.267
10	90	1.264	1.264	.162	7.850	.208	7.968
10	100	1.264	1.264	.154	8.550	.209	8.679
10	110	1.264	1.264	.146	9.250	.210	9.389
10	120	1.264	1.264	.138	9.950	.211	10.090
10	130	1.264	1.264	.130	10.650	.212	10.791
10	140	1.264	1.264	.122	11.350	.213	11.492
10	150	1.264	1.264	.114	12.050	.214	12.193
10	160	1.264	1.264	.106	12.750	.215	12.894
10	170	1.264	1.264	.098	13.450	.216	13.595
10	180	1.264	1.264	.090	14.150	.217	14.296
10	190	1.264	1.264	.082	14.850	.218	14.997
10	200	1.264	1.264	.074	15.550	.219	15.698
10	210	1.264	1.264	.066	16.250	.220	16.399
10	220	1.264	1.264	.058	16.950	.221	17.099
10	230	1.264	1.264	.050	17.650	.222	17.799
10	240	1.264	1.264	.042	18.350	.223	18.499
10	250	1.264	1.264	.034	19.050	.224	19.199
10	260	1.264	1.264	.026	19.750	.225	19.899
10	270	1.264	1.264	.018	20.450	.226	20.599
10	280	1.264	1.264	.010	21.150	.227	21.299
10	290	1.264	1.264	.002	21.850	.228	21.999
10	300	1.264	1.264	-.064	22.550	.229	22.699
10	310	1.264	1.264	-.143	23.250	.230	23.399
10	320	1.264	1.264	-.222	23.950	.231	24.099
10	330	1.264	1.264	-.299	24.650	.232	24.799
10	340	1.264	1.264	-.376	25.350	.233	25.499
10	350	1.264	1.264	-.453	26.050	.234	26.199
10	360	1.264	1.264	-.530	26.750	.235	26.899
10	370	1.264	1.264	-.607	27.450	.236	27.599
10	380	1.264	1.264	-.684	28.150	.237	28.299
10	390	1.264	1.264	-.761	28.850	.238	28.999
10	400	1.264	1.264	-.838	29.550	.239	29.699
10	410	1.264	1.264	-.915	30.250	.240	30.399
10	420	1.264	1.264	-.992	30.950	.241	31.099
10	430	1.264	1.264	-.969	31.650	.242	31.799
10	440	1.264	1.264	-.946	32.350	.243	32.499
10	450	1.264	1.264	-.923	33.050	.244	33.199
10	460	1.264	1.264	-.900	33.750	.245	33.899
10	470	1.264	1.264	-.877	34.450	.246	34.599
10	480	1.264	1.264	-.854	35.150	.247	35.299
10	490	1.264	1.264	-.831	35.850	.248	35.999
10	500	1.264	1.264	-.808	36.550	.249	36.699
10	510	1.264	1.264	-.785	37.250	.250	37.399
10	520	1.264	1.264	-.762	37.950	.251	38.099
10	530	1.264	1.264	-.739	38.650	.252	38.799
10	540	1.264	1.264	-.716	39.350	.253	39.499
10	550	1.264	1.264	-.693	40.050	.254	40.199
10	560	1.264	1.264	-.670	40.750	.255	40.899
10	570	1.264	1.264	-.647	41.450	.256	41.599
10	580	1.264	1.264	-.624	42.150	.257	42.299
10	590	1.264	1.264	-.601	42.850	.258	42.999
10	600	1.264	1.264	-.578	43.550	.259	43.699
10	610	1.264	1.264	-.555	44.250	.260	44.399
10	620	1.264	1.264	-.532	44.950	.261	44.699
10	630	1.264	1.264	-.509	45.650	.262	45.399
10	640	1.264	1.264	-.486	46.350	.263	46.099
10	650	1.264	1.264	-.463	47.050	.264	46.799
10	660	1.264	1.264	-.440	47.750	.265	47.499
10	670	1.264	1.264	-.417	48.450	.266	48.199
10	680	1.264	1.264	-.394	49.150	.267	48.899
10	690	1.264	1.264	-.371	49.850	.268	49.599
10	700	1.264	1.264	-.348	50.550	.269	50.299
10	710	1.264	1.264	-.325	51.250	.270	50.999
10	720	1.264	1.264	-.302	51.950	.271	51.699
10	730	1.264	1.264	-.279	52.650	.272	52.399
10	740	1.264	1.264	-.256	53.350	.273	53.099
10	750	1.264	1.264	-.233	54.050	.274	53.799
10	760	1.264	1.264	-.210	54.750	.275	54.499
10	770	1.264	1.264	-.187	55.450	.276	55.199
10	780	1.264	1.264	-.164	56.150	.277	55.899
10	790	1.264	1.264	-.141	56.850	.278	56.599
10	800	1.264	1.264	-.118	57.550	.279	57.299
10	810	1.264	1.264	-.095	58.250	.280	57.999
10	820	1.264	1.264	-.072	58.950	.281	58.699
10	830	1.264	1.264	-.049	59.650	.282	59.399
10	840	1.264	1.264	-.026	60.350	.283	60.099
10	850	1.264	1.264	-.003	61.050	.284	60.799
10	860	1.264	1.264	.160	61.750	.285	61.499
10	870	1.264	1.264	.183	62.450	.286	62.199
10	880	1.264	1.264	.206	63.150	.287	62.899
10	890	1.264	1.264	.229	63.850	.288	63.599
10	900	1.264	1.264	.252	64.550	.289	64.299
10	910	1.264	1.264	.275	65.250	.290	64.999
10	920	1.264	1.264	.298	65.950	.291	65.699
10	930	1.264	1.264	.321	66.650	.292	66.399
10	940	1.264	1.264	.344	67.350	.293	67.099
10	950	1.264	1.264	.367	68.050	.294	67.799
10	960	1.264	1.264	.390	68.750	.295	68.499
10	970	1.264	1.264	.413	69.450	.296	69.199
10	980	1.264	1.264	.436	70.150	.297	69.899
10	990	1.264	1.264	.459	70.850	.298	70.599
10	1000	1.264	1.264	.482	71.550	.299	71.299

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18 REFERENCES (Indicate on reverse side if necessary and identify by block number) 1. In this dissertation, Subset Selection, Stadeweise Procedure, Counter-Ordering, Order-Selection, Logistic Regression, Isotonic Regression, Non-Informative Prior, Posterior Distribution, and Conjugate priors are considered.		
19 ABSTRACT (Indicate on reverse side if necessary and identify by block number) The problem of selection and ranking (ordering) problems in statistics is often concerned with testing the classical truth of homogeneity hypothesis. Then the test statistic is used to select the treatments where the experimenter is interested in comparing the effectiveness of various treatments or processes with the goal of selecting the best treatment within a population.		
Chapter 1 of this thesis considers the problem of selection and ranking of treatments in a population if it is better than a control under the assumption that		

an ordering prior. Here, by an ordering prior we mean that there exists a known simple or partial order relationship among the unknown parameters of the treatments (excluding the control). Three new selection procedures are proposed and studied. These procedures do meet the usual requirement that the probability of a correct selection is greater than or equal to a pre-specified number P^* . Two of the three procedures use the isotonic regression estimator sample means of the k treatments with respect to the given ordering prior. Tables which are necessary to carry out the selection procedure with each of the three approaches for the selection of unknown means of normal populations and of non-normal populations are given. Monte Carlo comparisons on the performance of the three procedures for the normal or gamma mean problem were carried out in various selected cases. The results of this study seem to indicate that the procedures based on isotonic estimators always have superior performance, especially when there are more than one bad population. (in comparison with the existing)

Chapter II deals with a new "Bayes- P^* " approach about the procedure of selecting a subset which contains the "best" of k populations, where by best we mean the (unknown) population with the largest unknown mean. The (non-randomized) Bayes- P^* rule refers to a rule with minimax risk in the class of (non-randomized) rules which satisfy the condition that the probability of the ability of selecting the best is at least equal to P^* . Given the priors of the unknown parameters, two "Bayes- P^* " subset selection procedure,

J^B and J_{NR}^B (randomized and non-randomized, respectively) under certain loss functions are obtained and compared with the classical subset selection procedure J^M . The comparisons of the performance of J^B with J_{NR}^B and J^M , from Monte Carlo studies, indicate that the procedure J^B has higher power efficiency and smaller expected size of the selected subset. The results also indicate that J^B is robust when the true distributions are not normal; it is also robust to some other symmetric distributions such as, the logistic, the double exponential, the tri-modal (Laplace) and the gross error model (the contaminated distribution).

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